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Certifications

WBENC: 237019

1752439743100-86536

DBE: VN 20657

HUB:

NCTRCA WFWB38444Y0909

NELAP Certifications

Lubbock:

T104704219-08-TX

El Paso:

T104704221-08-TX

Midland: T104704392-08-TX

LELAP-02003 Kansas E-10317 LELAP-02002

Analytical and Quality Control Report

Brad Davis Zia Engineering & Environmental 755 S. Telshor Blvd. Suite F-201 Las Cruces, NM, 88011

Report Date: September 11, 2009

Work Order:

9082815

Project Name: HELSTF Chromate Spill Groundwater

Enclosed are the Analytical Report and Quality Control Report for the following sample(s) submitted to TraceAnalysis, Inc.

			Date	Time	Date
\mathbf{Sample}	Description	Matrix	Taken	Taken	Received
208427	HLSF-0143-HMW-037-0809	water	2009-08-26	09:55	2009-08-27

These results represent only the samples received in the laboratory. The Quality Control Report is generated on a batch basis. All information contained in this report is for the analytical batch(es) in which your sample(s) were analyzed.

This report consists of a total of 79 pages and shall not be reproduced except in its entirety, without written approval of TraceAnalysis, Inc.

Notes:

For inorganic analyses, the term MQL should actually read PQL.

Standard Flags

 ${f U}\,$ - Not detected. The analyte is not detected above the SDL.

 ${f J}$ - Estimated. The analyte is positively identified and the value is approximated between the SDL and MQL.

 $\, B \,\,$ - The sample contains less than ten times the concentration found in the method blank.

 ${f JB}$ - The analyte is positively identified and the value is approximated between the SDL and MQL.

The sample contains less than ten times the concentration found in the method blank.

The result should be considered non-detect to the SDL.

Dr. Blair Leftwich, Director
Dr. Michael Abel, Project Manager

Case Narrative

Samples for project HELSTF Chromate Spill Groundwater were received by TraceAnalysis, Inc. on 2009-08-27 and assigned to work order 9082815. Samples for work order 9082815 were received intact without headspace and at a temperature of 12.0 deg. C.

Samples were analyzed for the following tests using their respective methods.

		Prep	Prep	QC	${ m Analysis}$
Test	Method	Batch	Date	Batch	\mathbf{Date}
Ag, Total	S 6010B	53951	2009-09-03 at 08:16	63221	2009-09-03 at 10:42
Alkalinity	SM 2320B	54095	2009-09-08 at 11:00	63379	2009-09-08 at $11:00$
Al, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
${ m Ammonia}$	SM 4500-NH3 B,C	53864	2009-08-31 at $10:45$	63111	2009-08-31 at $16:00$
As, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Ba, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Be, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Bromide (IC)	$\to 300.0$	53984	2009-08-31 at $13:04$	63249	2009-08-31 at $20:24$
Ca, Total	S 6010B	53951	2009-09-03 at $08:16$	63355	2009-09-09 at $08:18$
Cd, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Chloride (IC)	$\to 300.0$	53984	2009-08-31 at $13:04$	63249	2009-08-31 at $20:24$
Chromium, Hexavalent	SM 3500 -Cr B	53776	2009-08-27 at 09:23	63015	2009-08-27 at 09:23
Co, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Cr, Dissolved	S 6010B	54153	2009-09-11 at $08:25$	63461	2009-09-11 at $11:54$
Cr, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Cu, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Explosives (8330)	S 8330-C18	53897	2009-09-01 at $15:00$	63142	2009-09-01 at $17:42$
Fe, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Fluoride (IC)	E 300.0	53984	2009-08-31 at $13:04$	63249	2009-08-31 at $20:24$
Hg, Total	S 7470A	53821	2009-08-31 at $12:30$	63072	2009-08-31 at $13:57$
K, Total	S 6010B	53951	2009-09-03 at $08:16$	63355	2009-09-09 at $08:18$
Mg, Total	S 6010B	53951	2009-09-03 at $08:16$	63355	2009-09-09 at $08:18$
Mn, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Mo, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Na, Total	S 6010B	53951	2009-09-03 at $08:16$	63355	2009-09-09 at $08:18$
Ni, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Nitrate and Nitrite as N	SM 4500-NO3 E	54188	2009-09-11 at $10:24$	63485	2009-09-11 at $16:25$
O/G	$\to 1664$	54129	2009-09-08 at $09:15$	63411	2009-09-09 at 11:48
Pb, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
pH	SM 4500-H+	53841	2009-08-27 at $13:50$	63082	2009-08-27 at 13:50
P, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Sb, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Semivolatiles	S 8270C	53996	2009-09-02 at $15:00$	63261	2009-09-04 at $10:45$
Se, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
SO4 (IC)	$\to 300.0$	53984	2009-08-31 at $13:04$	63249	2009-08-31 at $20:24$
TDS	SM 2540C	54007	2009-09-01 at $15:25$	63272	2009-09-01 at $15:25$
TKN	$\to 351.3$	54023	2009-09-02 at 11:00	63291	2009-09-02 at $13:30$
Tl, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
TOC	SM 5310C	54065	2009-09-08 at $13:26$	63340	2009-09-08 at $13:29$
Total Cyanide	SM 4500-CN C,E	53840	2009-08-31 at $10:15$	63080	2009-08-31 at 15:15
TPH DRO	Mod. 8015B	53812	2009-08-28 at 15:00	63042	2009-08-28 at 20:00

		Prep	Prep	QC	${ m Analysis}$
Test	Method	Batch	Date	Batch	Date
TPH GRO	S~8015B	53796	2009-08-28 at $14:49$	63025	2009-08-28 at $14:49$
V, Total	S_{010B}	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Zn, Total	S_{6010B}	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$

Results for these samples are reported on a wet weight basis unless data package indicates otherwise.

A matrix spike (MS) and matrix spike duplicate (MSD) sample is chosen at random from each preparation batch. The MS and MSD will indicate if a site specific matrix problem is occurring, however, it may not pertain to the samples for work order 9082815 since the sample was chosen at random. Therefore, the validity of the analytical data reported has been determined by the laboratory control sample (LCS) and the method blank (MB). These quality control measures are performed with each preparation batch to ensure data integrity.

All other exceptions associated with this report have been footnoted on the appropriate analytical page to assist in general data comprehension. Please contact the laboratory directly if there are any questions regarding this project.

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HELSTF Chromate Spill Groundwater

Analytical Report

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analytical Method: Analysis: Ag, Total S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQLMethod Based Based Blank MQLMDL SDLParameter Result Result Result Units Dilution (Unadjusted) (Unadjusted) Flag Total Silver 0.00111 < 0.00111 < 0.00500< 0.00111mg/L0.00111 0.005

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Al, Total S 3010A Analysis: Analytical Method: S 6010B Prep Method: QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQLMethod Based Based Blank MQLMDLFlag Parameter Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Aluminum 0.1350.135< 0.00301 mg/L0.003010.050.00301

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: El Paso

Analysis: Alkalinity Analytical Method: SM 2320BPrep Method: N/AQC Batch: 63379 Date Analyzed: 2009-09-08 Analyzed By: JGPrep Batch: Sample Preparation: Prepared By: 54095JG

SDLMQLMethod Based Based Blank MQLMDL Flag Parameter Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Hydroxide Alkalinity < 1.00 <1.00 < 1.00 mg/L as CaCo3 1.00 1 1 UCarbonate Alkalinity < 1.00 < 1.00 < 1.00mg/L as CaCo3 1 1.00 1 1 Bicarbonate Alkalinity 122122 < 4.00 mg/L as CaCo3 4.00 4 4 1 mg/L as CaCo3Total Alkalinity 122 122 < 4.001 4.00 4 4

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Analytical Method: SM 4500-NH3 B,C Prep Method: N/AAmmonia QC Batch: Date Analyzed: Analyzed By: AH63111 2009-08-31 Prep Batch: 53864Sample Preparation: 2009-08-31 Prepared By: AH

Report Date: September 11, 2009 Work Order: 9082815 Page Number: 6 of 79

HELSTF Chromate Spill Groundwater

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Ammonia-N	J	0.392	<1.00	< 0.353	mg/L	1	0.353	1	0.353

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: As, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 2009-09-03 Analyzed By: 63221 Date Analyzed: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQL Method Based Based Blank MQLMDL Parameter Result Result Result Units Dilution SDL Flag (Unadjusted) (Unadjusted) Total Arsenic < 0.00448 < 0.0100 < 0.00448mg/L0.00448 0.01 0.00448 1

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Ba, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: Analyzed By: 2009-09-03 RRPrep Batch: Sample Preparation: 2009-09-03 Prepared By: KV53951

SDLMQLMethod MQLMDL Based Based Blank Parameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Barium 0.01300.0130< 0.00105 0.001050.005 0.00105 mg/L1

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Be, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: Date Analyzed: 2009-09-03 Analyzed By: RR63221 Prep Batch: 53951Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQLMethod MQLMDL Based Based Blank Parameter Flag Result Result Result Dilution SDL(Unadjusted) Units (Unadjusted) Total Beryllium < 0.000450 < 0.00200 < 0.000450 mg/L0.0004500.002 0.00045

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Bromide (IC) Analytical Method: E 300.0 Prep Method: N/A

Report Date	e: Septem	ber 11, 200	9		Vork Order: Chromate S	9082815 Spill Ground	water	Page Number: 7 of 79		
QC Batch: Prep Batch:			Date Anal Sample Pr	yzed: eparation:	2009-08-31 2009-08-31		Analyz Prepar	•		
D	TO!	$\begin{array}{c} \mathrm{SDL} \\ \mathrm{Based} \end{array}$	MQL Based	Method Blank	TT **	Dil et	aDI.	MQL	MDL	
Parameter	Flag	Result	Result	Result	Units	$\operatorname{Dilution}$	SDL	(Unadjusted)	(Unadjusted)	
$\overline{\text{Bromide}}$	U	< 0.960	< 2.50	< 0.960	${ m mg/L}$	5	0.960	0.5	0.192	

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory:	$\operatorname{Lubbock}$								
Analysis:	Ca, Total		Ana	lytical Meth	od: S	6010B		Prep Meth	od: S 3010A
QC Batch:	63355		Date	e Analyzed:	20	009-09-09		Analyzed E	By: RR
Prep Batch:	53951		Sam	ple Preparat	ion: 20	009-09-03		Prepared E	By: KV
		SDL	MQL	Method					
		Based	Based	Blank				$_{ m MQL}$	MDL
Parameter	Flag	Result	Result	Result	Units	$\operatorname{Dilution}$	SDL	(Unadjusted)	(Unadjusted)
Total Calciur	n	544	544	<1.17	mg/L	10	1.17	1	0.117

Sample: 208427 - HLSF-0143-HMW-037-0809

Analysis: Co QC Batch: 63	Cd, Total 63221		Date A	ical Method: malyzed: Preparation	2009-	10B -09-03 -09-03		Prep Metho Analyzed B Prepared B	·
		$_{ m Based}^{ m SDL}$	$egin{array}{l} \mathrm{MQL} \\ \mathrm{Based} \end{array}$	$egin{array}{c} ext{Method} \ ext{Blank} \end{array}$				$_{ m MQL}$	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Cadmium	U	< 0.000303	< 0.00200	< 0.000303	mg/L	1	0.000303	0.002	0.000303

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory:	Lubboc	k								
Analysis:	Chlorid	e (IC)		Analytical	Method:	E 300.0		Prep Method: N/A		
QC Batch:	63249			Date Anal	lyzed:	2009-08-31		Analyz	ed By: SS	
Prep Batch:	53984			Sample Pr	reparation:	2009-08-31		Prepar	ed By: SS	
		SDL	MQL	Method						
		Based	Based	Blank				MQL	MDL	
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)	
Chloride		220	220	< 0.785	$\mathrm{mg/L}$	5	0.785	2.5	0.157	

Sample: 208427 - HLSF-0143-HMW-037-0809

Report Date: September 11, 2009 Work Order: 9082815 Page Number: 8 of 79

HELSTF Chromate Spill Groundwater

Laboratory: El Paso

Analysis: Analytical Method: Chromium, Hexavalent SM 3500-Cr B Prep Method: N/AQC Batch: Date Analyzed: 2009-08-27 Analyzed By: MD63015 Prep Batch: 53776 Sample Preparation: 2009-08-27 Prepared By: MD

> SDLMQLMethod

Based Based Blank MQLMDL Parameter Flag Result Result ResultUnits Dilution SDL(Unadjusted) (Unadjusted) Hexavalent Chromium < 0.00594< 0.0100 < 0.00594 mg/L0.005940.01 0.00594

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Co, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQL Method Based Based Blank

MQLMDLParameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Cobalt < 0.000822 < 0.00200 < 0.000822 mg/L0.0008220.002 0.000822

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Analytical Method: Prep Method: S 3005A Cr, Dissolved S 6010B QC Batch: 63461 Date Analyzed: 2009-09-11 Analyzed By: RRPrep Batch: 54153Sample Preparation: 2009-09-11 Prepared By: KV

SDLMQL Method Based Based Blank

MQLMDL SDLParameter Flag Result Result Result Units Dilution (Unadjusted) (Unadjusted) Dissolved Chromium 0.003000.003000.0010.000583 < 0.000583 mg/L0.000583

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Cr, Total Analytical Method: Prep Method: S 3010A S 6010B QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951Sample Preparation: 2009-09-03 Prepared By: KV

SDL MQLMethod

		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Chromium	J	0.00400	< 0.00500	< 0.000583	mg/L	1	0.000583	0.005	0.000583

Report Date: September 11, 2009 Work Order: 9082815 Page Number: 9 of 79

HELSTF Chromate Spill Groundwater

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Cu, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

 SDL MQL Method

Blank MQLBased Based MDL Dilution Parameter Flag Result Result Result Units SDL(Unadjusted) (Unadjusted) Total Copper < 0.000843 < 0.00500 < 0.000843 mg/L 0.0008430.0050.000843

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Explosives (8330) Analytical Method: S 8330-C18 Prep Method: S 3535A QC Batch: Date Analyzed: 2009-09-01 Analyzed By: DS63142Sample Preparation: Prep Batch: 538972009-09-01 Prepared By: DS

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	$\operatorname{Dilution}$	SDL	(Unadjusted)	(Unadjusted)
HMX	U	< 0.123	< 0.500	< 0.123	$\mu \mathrm{g/L}$	1	0.123	0.5	0.123
RDX	U	< 0.298	< 0.500	< 0.298	$\mu { m g/L}$	1	0.298	0.5	0.298
1,3,5-Trinitrobenzene	U	< 0.339	< 0.500	< 0.339	$\mu { m g/L}$	1	0.339	0.5	0.339
1,3-Dinitrobenzene	U	< 0.389	< 0.500	< 0.389	$\mu { m g/L}$	1	0.389	0.5	0.389
Nitrobenzene	U	< 0.379	< 0.500	< 0.379	$\mu { m g/L}$	1	0.379	0.5	0.379
Tetryl	U	< 0.413	< 0.500	< 0.413	$\mu { m g/L}$	1	0.413	0.5	0.413
TNT	U	< 0.464	< 0.500	< 0.464	$\mu { m g/L}$	1	0.464	0.5	0.464
4-Amino-DNT	U	< 0.319	< 0.500	< 0.319	$\mu { m g/L}$	1	0.319	0.5	0.319
2-Amino-DNT	U	< 0.391	< 0.500	< 0.391	$\mu { m g/L}$	1	0.391	0.5	0.391
2,6-DNT	U	< 0.323	< 0.500	< 0.323	$\mu \mathrm{g/L}$	1	0.323	0.5	0.323
2,4-DNT	U	< 0.366	< 0.500	< 0.366	$\mu { m g/L}$	1	0.366	0.5	0.366
2-NT	U	< 0.379	< 0.500	< 0.379	$\mu { m g/L}$	1	0.379	0.5	0.379
4-NT	U	< 0.398	< 0.500	< 0.398	$\mu { m g/L}$	1	0.398	0.5	0.398
3-NT	U	< 0.346	< 0.500	< 0.346	μ_g/L	1	0.346	0.5	0.346

					Spike	$\operatorname{Percent}$	$\operatorname{Recovery}$
$\mathbf{Surrogate}$	Flag	Result	${ m Units}$	Dilution	${f Amount}$	Recovery	Limits
1,2-Dinitrobenzene		1.87	$\mu \mathrm{g/L}$	1	2.50	75	19.8 - 160

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Fe, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951Sample Preparation: 2009-09-03 Prepared By: KV

Report Date: September 11, 2009 Work Order: 9082815 Page Number: 10 of 79

HELSTF Chromate Spill Groundwater

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	$\operatorname{Dilution}$	SDL	(Unadjusted)	(Unadjusted)
Total Iron		0.0360	0.0360	< 0.000872	mg/L	1	0.000872	0.01	0.000872

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Fluoride (IC) Analytical Method: E 300.0 Prep Method: N/AQC Batch: 63249 Analyzed By: Date Analyzed: 2009-08-31 SSPrep Batch: 53984 Sample Preparation: 2009-08-31 Prepared By: SS

SDLMQLMethod Based Based Blank MQLMDL Parameter Result Units Dilution SDLFlag Result Result (Unadjusted) (Unadjusted) Fluoride <1.02 < 2.50<1.02 mg/L1.02 0.5 0.204 5

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

N/AAnalysis: Hg, Total Analytical Method: S 7470A Prep Method: QC Batch: 63072 Analyzed By: TPDate Analyzed: 2009-08-31 Prep Batch: Sample Preparation: 2009-08-31 Prepared By: TP53821

SDLMQL Method Based MQLMDL Based Blank Parameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Mercury < 0.0000329 < 0.000200 < 0.0000329 0.00003290.0002 3.29e-05mg/L

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: K, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RRPrep Batch: 53951Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQLMethod MQLMDL Based Based Blank Parameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Potassium 86.2 86.2 < 0.172mg/L0.172 0.172

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Mg, Total Analytical Method: S 6010B Prep Method: S 3010A

Report Date: September 11, 2009 Work Order: 9082815 Page Number: 11 of 79 HELSTF Chromate Spill Groundwater QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RRPrep Batch: Sample Preparation: 2009-09-03 53951Prepared By: KVSDLMQLMethod Based Blank MQLBased MDLParameter Flag ${\bf Result}$ Result Result Units Dilution SDL(Unadjusted) (Unadjusted)

mg/L

10

1.60

0.16

< 1.60

Sample: 208427 - HLSF-0143-HMW-037-0809

328

328

Total Magnesium

Laboratory:	${ m Lubbock}$								
Analysis:	Mn, Total		Analyt	ical Method:	S 601	.0B		Prep Metho	d: S 3010A
QC Batch:	63221		$\operatorname{Date} A$	Date Analyzed:		09-03		Analyzed B	y: RR
Prep Batch:	53951		$_{ m Sample}$	Preparation	: 2009-	09-03		Prepared B	y: KV
		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Manga	nese J	0.000400	< 0.00250	< 0.000305	mg/L	1	0.000305	0.0025	0.000305

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory:	Lubbock									
Analysis:	Mo, Tota	$_{ m al}$		Analy	tical Metho	d: S 6	010B		Prep Metho	od: S 3010A
QC Batch:	63221			Date	Analyzed:	200	9-09-03		Analyzed B	y: RR
Prep Batch:	53951			Samp	le Preparati	on: 200	9-09-03		Prepared B	y: KV
			SDL	MQL	Method					
			Based	Based	Blank				MQL	MDL
Parameter]	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Molybd	lenum		0.125	0.125	< 0.00119	mg/L	1	0.00119	0.01	0.00119

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Analysis: QC Batch: Prep Batch:	Lubbock Na, Total 63355 53951		Dat	alytical Metl se Analyzed: aple Prepara	2	6 6010B 1009-09-09 1009-09-03		Prep Meth Analyzed E Prepared E	By: RR
		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	$\operatorname{Dilution}$	SDL	(Unadjusted)	(Unadjusted)
Total Sodium	n	401	401	< 0.500	mg/L	10	0.500	1	0.05

Sample: 208427 - HLSF-0143-HMW-037-0809

Report Date: September 11, 2009 Work Order: 9082815 Page Number: 12 of 79

HELSTF Chromate Spill Groundwater

Laboratory: Lubbock Analysis: Ni, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: Date Analyzed: 2009-09-03 Analyzed By: 63221 RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KVSDLMQL Method Based Based Blank MQLMDL

Parameter Flag ${\bf Result}$ Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Nickel < 0.00121 < 0.00500< 0.00121mg/L0.00121 0.0050.00121

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Nitrate and Nitrite as N Analytical Method: SM 4500-NO3 E Prep Method: N/AQC Batch: 63485 Date Analyzed: 2009-09-11 Analyzed By: KVPrep Batch: 54188Sample Preparation: 2009-09-11 Prepared By: KV

SDL MQLMethod Based MQLMDLBased Blank Parameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) 7.00 Nitrate and Nitrite as N 14.5<20.0 < 7.00mg/L200 0.1 0.035

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: El Paso

O/G Analytical Method: Analysis: E 1664 Prep Method: N/AQC Batch: 63411 Date Analyzed: 2009-09-09 Analyzed By: MDPrep Batch: 54129Sample Preparation: 2009-09-08 Prepared By: MD

SDLMQL Method Based Based Blank MQLMDL Flag SDL Parameter Result Result Result Units Dilution (Unadjusted) (Unadjusted) Oil and Grease < 3.60< 5.00< 3.60mg/L 3.60 5 3.6

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: P, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDLMethod MQL Based Based Blank MQLMDLParameter Result Result Result Units Dilution SDLFlag (Unadjusted) (Unadjusted) Total Phosphorous < 0.00289 < 0.0250< 0.00289 mg/L0.002890.0250.00289

Report Date: September 11, 2009 Work Order: 9082815 Page Number: 13 of 79

HELSTF Chromate Spill Groundwater

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Pb, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

 SDL MQL Method

Blank MQLBased Based MDL Parameter Result Result Result Units Dilution SDL(Unadjusted) Flag (Unadjusted) Total Lead < 0.00326 < 0.00500 < 0.00326 mg/L 0.00326 0.0050.00326

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: El Paso

Analysis: рΗ Analytical Method: SM 4500-H+Prep Method: N/AQC Batch: 63082 JGDate Analyzed: 2009-08-27 Analyzed By: Prep Batch: 53841 Sample Preparation: 2009-08-27 Prepared By: JR

RL

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Sb. Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDL MQL Method Based Based Blank

Based Based Blank MQLMDL Flag SDLParameter Result Result Result Units Dilution (Unadjusted) (Unadjusted) 0.02 Total Antimony < 0.00440< 0.0200< 0.00440mg/L0.004400.0044

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Se, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951KVSample Preparation: 2009-09-03 Prepared By:

 $\mathrm{SDL} \qquad \mathrm{MQL} \qquad \mathrm{Method}$

		Basea	Basea	Biank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Selenium	J	0.0140	< 0.0200	< 0.00508	$\mathrm{mg/L}$	1	0.00508	0.02	0.00508

MAT

MINI

HELSTF Chromate Spill Groundwater

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory:	Lubbock				
Analysis:	Semivolatiles	Analytical Method:	S 8270 C	Prep Method:	S 3510 C
QC Batch:	63261	Date Analyzed:	2009-09-04	Analyzed By:	MN
Prep Batch:	53996	Sample Preparation:	2009-09-02	Prepared By:	MN

		SDL	MQL	Method				
		Based	Based	Blank			MQL	MDL
Parameter	Flag	Result	Result	Result Ur	its Dilution	$_{ m l}$ SDL	(Unadjusted)	(Unadjusted)
Pyridine	U	< 0.000624	< 0.00513	<0.000624 mg	/L 1.026	0.000624	0.005	0.000608
N-Nitrosodimethylamine	U	< 0.000566	< 0.00513	<0.000566 mg	/L 1.026	0.000566	0.005	0.000552
2-Picoline	U	< 0.000419	< 0.00513	<0.000419 mg	/L 1.026	0.000419	0.005	0.000408
Methyl methanesulfonate	U	< 0.000359	< 0.00513	<0.000359 mg	/L 1.026	0.000359	0.005	0.00035
Ethyl methanesulfonate	U	< 0.000460	< 0.00513	<0.000460 mg	/L 1.026	0.000460	0.005	0.000448
Phenol	U	< 0.000522	< 0.00513	<0.000522 mg	/L 1.026	0.000522	0.005	0.000509
Aniline	U	< 0.000709	< 0.00513	<0.000709 mg	/L 1.026	0.000709	0.005	0.000691
${ m bis}(2{ ext{-chloroet}}{ m hyl}){ m et}{ m her}$	U	< 0.000451	< 0.00513	<0.000451 mg	/L 1.026	0.000451	0.005	0.00044
2-Chlorophenol	U	< 0.000551	< 0.00513	< 0.000551 mg	L/L 1.026	0.000551	0.005	0.000537
1,3-Dichlorobenzene (meta)	U	< 0.000452	< 0.00513	<0.000452 mg	L/L 1.026	0.000452	0.005	0.000441
1,4-Dichlorobenzene (para)	U	< 0.000451	< 0.00513	<0.000451 mg	L/L 1.026	0.000451	0.005	0.00044
Benzyl alcohol	U	< 0.000552	< 0.00513	< 0.000552 mg	L/L 1.026	0.000552	0.005	0.000538
1,2-Dichlorobenzene (ortho)	U	< 0.000454	< 0.00513	<0.000454 mg	L/L 1.026	0.000454	0.005	0.000443
2-Methylphenol	U	$< 0.000745 \cdot$	< 0.00513	<0.000745 mg	L/L 1.026	0.000745	0.005	0.000726
${ m bis}(2{ ext{-chloroisopropyl}}) { m ether}$	U	< 0.000516	< 0.00513	<0.000516 mg	L/L 1.026	0.000516	0.005	0.000503
4-Methylphenol / 3-Methylphenol		$< 0.000525 \cdot$	< 0.00513	< 0.000525 mg	L/L 1.026	0.000525	0.005	0.000512
N-Nitrosodi-n-propylamine	U	< 0.000751	< 0.00513	<0.000751 mg	L/L 1.026	0.000751	0.005	0.000732
$\operatorname{Hexachloroethane}$	U	< 0.000520	< 0.00513	<0.000520 mg	L/L 1.026	0.000520	0.005	0.000507
$egin{array}{c} Acetophenone \end{array}$	U	$< 0.000435 \cdot$	< 0.00513	<0.000435 mg	L/L 1.026	0.000435	0.005	0.000424
Nitrobenzene	U	$< 0.000477 \cdot$	< 0.00513	$< 0.000477 \mathrm{mg}$	L/L 1.026	0.000477	0.005	0.000465
N-Nitrosopiperidine	U	< 0.000454	< 0.00513	<0.000454 mg	L/L 1.026	0.000454	0.005	0.000443
Isophorone	U	< 0.000635	< 0.00513	< 0.000635 mg	L/L = 1.026	0.000635	0.005	0.000619
2-Nitrophenol	U	< 0.000416	< 0.00513	< 0.000416 mg	L/L = 1.026	0.000416		0.000406
2,4-Dimethylphenol	U	$< 0.000489 \cdot$	< 0.00513	<0.000489 mg	L/L 1.026	0.000489	0.005	0.000477
${ m bis}(2{ m -chloroethoxy}){ m methane}$	U	< 0.000443	< 0.00513	<0.000443 mg		0.000443	0.005	0.000432
2,4-Dichlorophenol	U	< 0.000410	< 0.00513	<0.000410 mg	L/L = 1.026	0.000410	0.005	0.0004
1,2,4-Trichlorobenzene	U			<0.000414 mg		0.000414	0.005	0.000404
Benzoic acid	U	< 0.00167	< 0.00513	$< 0.00167 \mathrm{mg}$	L/L 1.026	0.00167	0.005	0.00163
Naphthalene	U	< 0.000502	< 0.00513	<0.000502 mg	L/L 1.026	0.000502	0.005	0.000489
a,a-Dimethylphenethylamine	U	< 0.00132	< 0.00513	< 0.00132 mg	L/L 1.026	0.00132	0.005	0.00129
4-Chloroaniline	U	< 0.000388	< 0.00513	< 0.000388 mg	L/L = 1.026	0.000388	0.005	0.000378
2,6-Dichlorophenol	U	< 0.000496	< 0.0103	< 0.000496 mg	L/L = 1.026	0.000496	0.01	0.000484
${\it Hexachlorobut}$ adiene	U	< 0.000530	< 0.00513	< 0.000530 mg	L/L = 1.026	0.000530	0.005	0.000517
N-Nitroso-di-n-butylamine	U	< 0.000673	< 0.00513	< 0.000673 mg	L/L = 1.026	0.000673	0.005	0.000656
4-Chloro-3-methylphenol	U	< 0.000536	< 0.00513	< 0.000536 mg	L/L = 1.026	0.000536	0.005	0.000522
$2 ext{-Methylnaphthalene}$	U	< 0.000434	< 0.00513	<0.000434 mg	L/L 1.026	0.000434	0.005	0.000423
$1 ext{-} ext{Methylnaphthalene}$	U	< 0.000508	< 0.00513	< 0.000508 mg	L/L = 1.026	0.000508	0.005	0.000495
1,2,4,5-Tetrachlorobenzene	U	< 0.000628	< 0.00513	<0.000628 mg	L/L = 1.026	0.000628	0.005	0.000612
${\it Hexachlorocyclopenta}$ diene	U	< 0.000572	< 0.00513	<0.000572 mg		0.000572	0.005	0.000558
2,4,6-Trichlorophenol	U	< 0.000815	< 0.0103	< 0.000815 mg	/L 1.026	0.000815	0.01	0.000794

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sample 208427 continued . . .

		SDL	MQL	Method				
		Based	Based	Blank			MQL	MDL
Parameter	Flag	Result	Result	Result Units	Dilution	$_{ m L}$ SDL	(Unadjusted) (Unadjusted)
$\overline{2,4,5}$ -Trichlorophenol	U			$< 0.000856 \mathrm{mg/L}$	1.026	0.000856	0.005	0.000834
$2 ext{-Chloronaphthalene}$	U	< 0.000427 <	< 0.00513	$< 0.000427 \mathrm{mg/L}$	1.026	0.000427	0.005	0.000416
$1 ext{-Chloronaphthalene}$	U	< 0.000488	< 0.00513	$< 0.000488 \mathrm{mg/L}$	1.026	0.000488	0.005	0.000476
2-Nitroaniline	U	< 0.000780	< 0.00513	< 0.000780 mg/L	1.026	0.000780	0.005	0.00076
${\bf Dimethylphthalate}$	U	< 0.000660	< 0.00513	$< 0.000660 \mathrm{mg/L}$	1.026	0.000660	0.005	0.000643
${ m Acenaphthylene}$	U	< 0.000601	< 0.00513	< 0.000601 mg/L	1.026	0.000601	0.005	0.000586
2,6-Dinitrotoluene	U	< 0.000657 <	< 0.00513	$< 0.000657 \mathrm{mg/L}$	1.026	0.000657	0.005	0.00064
3-Nitroaniline	U	< 0.000740	< 0.00513	< 0.000740 mg/L	1.026	0.000740	0.005	0.000721
${ m Acenaphthene}$	U	< 0.000434	< 0.00513	< 0.000434 mg/L	1.026	0.000434	0.005	0.000423
2,4-Dinitrophenol	U	< 0.000226	< 0.00513	$< 0.000226 \mathrm{mg/L}$	1.026	0.000226	0.005	0.00022
Dibenzofuran	U	< 0.000419	< 0.00513	$< 0.000419 \mathrm{mg/L}$	1.026	0.000419	0.005	0.000408
Pentachlorobenzene	U	< 0.000586	< 0.00513	$< 0.000586 \mathrm{mg/L}$	1.026	0.000586	0.005	0.000571
4-Nitrophenol	U	< 0.00190	< 0.0256	$< 0.00190 \mathrm{mg/L}$	1.026	0.00190	0.025	0.00185
2,4-Dinitrotoluene	J	0.00103 <	< 0.00513	< 0.000935 mg/L	1.026	0.000935	0.005	0.000911
1-Naphthylamine	U	< 0.000706	< 0.00513	$< 0.000706 \mathrm{mg/L}$	1.026	0.000706	0.005	0.000688
2,3,4,6-Tetrachlorophenol	U			$< 0.000580 \mathrm{mg/L}$	1.026	0.000580	0.01	0.000565
2-Naphthylamine	U			$< 0.000717 \mathrm{mg/L}$	1.026	0.000717	0.005	0.000699
Fluorene	U			< 0.000665 mg/L	1.026	0.000665	0.005	0.000648
4-Chlorophenyl-phenylether	U			< 0.000635 mg/L	1.026	0.000635	0.005	0.000619
Diethylphthalate	U			$< 0.000850 \mathrm{mg/L}$	1.026	0.000850	0.005	0.000828
4-Nitroaniline	U			$< 0.000720 \mathrm{mg/L}$	1.026	0.000720	0.005	0.000702
Diphenylhydrazine	U			$< 0.000674 \mathrm{mg/L}$	1.026	0.000674	0.005	0.000657
4,6-Dinitro-2-methylphenol	U			$< 0.00203 \mathrm{mg/L}$	1.026	0.00203	0.005	0.00198
Diphenylamine	U			< 0.000451 mg/L	1.026	0.000451	0.005	0.00044
4-Bromophenyl-phenylether	U			$< 0.000564 \mathrm{mg/L}$	1.026	0.000564	0.005	0.00055
Phenacetin	U			< 0.000621 mg/L	1.026	0.000621	0.005	0.000605
Hexachlorobenzene	U			$< 0.000519 \mathrm{mg/L}$	1.026	0.000519	0.005	0.000506
4-Aminobiphenyl	U			< 0.000541 mg/L	1.026	0.000541	0.005	0.000527
Pentachlorophenol	U			$< 0.000446 \mathrm{mg/L}$	1.026	0.000446	0.01	0.000435
Anthracene	U			$< 0.000439 \mathrm{mg/L}$	1.026	0.000439	0.005	0.000428
Pentachloronitrobenzene	U			$< 0.000419 \mathrm{mg/L}$	1.026	0.000419	0.005	0.000408
Pronamide	U			$< 0.000488 \mathrm{mg/L}$	1.026	0.000488	0.005	0.000476
Phenanthrene	U			< 0.000562 mg/L	1.026	0.000562	0.005	0.000548
Di-n-butylphthalate	U			$< 0.000496 \mathrm{mg/L}$	1.026	0.000496	0.005	0.000483
Fluoranthene	U			$< 0.000648 \mathrm{mg/L}$	1.026	0.000648	0.005	0.000632
Benzidine	U			< 0.00244 mg/L	1.026	0.00244	0.025	0.00238
Pyrene	U			< 0.000742 mg/L	1.026	0.000742	0.005	0.000723
p-Dimethylaminoazobenzene	U			< 0.000925 mg/L	1.026	0.000925	0.005	0.000902
Butylbenzylphthalate	U			$< 0.000456 \mathrm{mg/L}$	1.026	0.000456	0.005	0.000445
Benzo(a)anthracene	U			< 0.000541 mg/L	1.026	0.000541	0.005	0.000527
3,3-Dichlorobenzidine	U			< 0.00121 mg/L	1.026	0.00121	0.005	0.00118
Chrysene	U			< 0.000654 mg/L	1.026	0.000654	0.005	0.000638
bis(2-ethylhexyl)phthalate	U			$< 0.000576 \mathrm{mg/L}$	1.026	0.000576	0.005	0.000561
Di-n-octylphthalate	U			$< 0.00119 \mathrm{mg/L}$	1.026	0.00119	0.005	0.00116
Benzo(b)fluoranthene	U			< 0.000902 mg/L		0.000902	0.005	0.000879

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HELSTF Chromate Spill Groundwater

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sample 208427 continued . . .

		SDL	MQL	Method				
		Based	Based	Blank			MQL	MDL
Parameter	Flag	Result	Result	Result Units	Dilution	SDL	(Unadjusted)) (Unadjusted)
$\overline{\mathrm{Benzo}(\mathtt{k})}$ fluoranthene	U	< 0.000867	< 0.00513	$< 0.000867 \mathrm{mg/L}$	1.026	0.000867	0.005	0.000845
7,12-Dimethylbenz(a)anthracene	U	< 0.00105	< 0.00513	< 0.00105 mg/L	1.026	0.00105	0.005	0.00102
$\mathrm{Benzo}(\mathrm{a})\mathrm{pyrene}$	U	< 0.00171	< 0.00513	< 0.00171 mg/L	1.026	0.00171	0.005	0.00167
3-Methylcholanthrene	U	< 0.000932	< 0.00513	${<}0.000932~\rm{mg/L}$	1.026	0.000932	0.005	0.000908
${ m Dibenzo(a,j)}$ acridine	U	< 0.00132	< 0.00513	< 0.00132 mg/L	1.026	0.00132	0.005	0.00129
Indeno(1,2,3-cd)pyrene	U	< 0.000884	< 0.00513	${<}0.000884\:\mathrm{mg/L}$	1.026	0.000884	0.005	0.000862
${ m Dibenzo(a,h)}$ anthracene	U	< 0.000830	< 0.00513	< 0.000830 mg/L	1.026	0.000830	0.005	0.000809
$\mathrm{Benzo}(\mathrm{g,h,i})$ perylene	U	< 0.000974	< 0.00513	${<}0.000974\:\mathrm{mg/L}$	1.026	0.000974	0.005	0.000949

					Spike	$\operatorname{Percent}$	Recovery
Surrogate	Flag	Result	Units	Dilution	${f Amount}$	Recovery	Limits
2-Fluorophenol		0.0237	$_{ m mg/L}$	1.026	0.0800	30	10 - 53.1
${ m Phenol-d5}$		0.0157	${ m mg/L}$	1.026	0.0800	20	10 - 36.9
${ m Nitrobenzene-d5}$		0.0356	${ m mg/L}$	1.026	0.0800	44	23.8 - 108
2-Fluorobiphenyl		0.0248	$\mathrm{mg/L}$	1.026	0.0800	31	15.9 - 127
2,4,6-Tribromophenol		0.0539	$\mathrm{mg/L}$	1.026	0.0800	67	10 - 123
Terphenyl-d14		0.0439	m mg/L	1.026	0.0800	55	17.2 - 160

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: SO4 (IC) Analytical Method: E 300.0 Prep Method: N/AQC Batch: Date Analyzed: 2009-08-31 Analyzed By: SS 63249Prep Batch: 53984 Sample Preparation: 2009-08-31 Prepared By: SS

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
$\overline{\mathrm{Sulfate}}$		3090	3090	<15.3	$\mathrm{mg/L}$	100	15.3	2.5	0.153

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: El Paso

Analysis: TDSAnalytical Method: $\rm SM~2540C$ Prep Method: N/A QC Batch: 63272Date Analyzed: Analyzed By: MD2009-09-01 Prep Batch: 54007 Sample Preparation: 2009-09-01 Prepared By: MD

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Dissolved Solids		5320	5320	< 5.00	mg/L	1	5.00		5

HELSTF Chromate Spill Groundwater

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: TKN Analytical Method: E 351.3 Prep Method: N/AQC Batch: 63291 Date Analyzed: 2009-09-02 Analyzed By: AHPrep Batch: 54023 Sample Preparation: 2009-09-02 Prepared By: AH

SDL MQL Method

MQLMDLBased Based Blank Parameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Kjeldahl Nitrogen - N 3.64 <10.0 < 2.45mg/L 2.4510 2.45 1

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Tl, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 2009-09-03 Analyzed By: 63221 Date Analyzed: RRPrep Batch: 53951Sample Preparation: 2009-09-03 Prepared By: KV

SDL MQL Method
Based Blank

MQLMDL Based Based Blank SDLDilution Parameter Flag Result Result Result Units (Unadjusted) (Unadjusted) Total Thallium < 0.00488< 0.0500 < 0.00488mg/L1 0.00488 0.050.00488

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: TOC Analytical Method: SM 5310CPrep Method: N/AQC Batch: 63340 Date Analyzed: 2009-09-08 Analyzed By: KVPrep Batch: 54065Sample Preparation: 2009-09-08 Prepared By: KV

 $\mathrm{SDL} \quad \mathrm{MQL} \quad \mathrm{Method}$

Based MQLMDL Based Blank (Unadjusted) Parameter Result Result Result Dilution SDLFlag Units (Unadjusted) Total Organic Carbon 1.441.44< 0.401mg/L0.4010.401

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Total Cyanide Analytical Method: SM 4500-CN C,E Prep Method: N/AQC Batch: 63080 2009-08-31 AHDate Analyzed: Analyzed By: Prep Batch: 53840 Sample Preparation: 2009-08-31 Prepared By: AH

 $continued \dots$

HELSTF Chromate Spill Groundwater

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sample 208427 continued ...

Parameter	Flag	$\begin{array}{c} \mathrm{SDL} \\ \mathrm{Based} \\ \mathrm{Result} \end{array}$	$egin{array}{l} \mathrm{MQL} \\ \mathrm{Based} \\ \mathrm{Result} \end{array}$	Method Blank Result	$_{ m Units}$	Dilution	SDL	m MQL (Unadjusted)	MDL (Unadjusted)
		SDL	$_{ m MQL}$	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Cyanide	U	< 0.0110	< 0.0150	< 0.0110	mg/L	1	0.0110	0.015	0.011

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: TPH DRO Analytical Method: Prep Method: N/A Mod. 8015B

QC Batch: 63042 Date Analyzed: 2009-08-28 Analyzed By: Prep Batch: Sample Preparation: Prepared By: 538122009-08-28

SDLMQLMethod Based Based Blank MQLMDLSDLParameter Flag Result Result Result Units Dilution (Unadjusted) (Unadjusted) DRO < 0.8760.876 < 0.876< 5.00mg/L5 0.876

					Spike	Percent	Recovery
$\mathbf{Surrogate}$	Flag	Result	Units	Dilution	${f Amount}$	Recovery	Limits
n-Triacontane		8.39	$_{ m mg/L}$	1	10.0	84	34.4 - 185

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: TPH GRO Analytical Method: S 8015BPrep Method: S 5030B QC Batch: 63025 Date Analyzed: 2009-08-28 Analyzed By: ERPrep Batch: Sample Preparation: 2009-08-28 53796 Prepared By: ER

SDLMQLMethod Based Based Blank MQLMDLParameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) GRO < 0.0702 < 0.200 0.0702 < 0.0702mg/L0.07020.2 1

Surrogate	Flag	Result	Units	Dilution	$\begin{array}{c} {\rm Spike} \\ {\rm Amount} \end{array}$	Percent Recovery	$rac{ m Recovery}{ m Limits}$
Trifluorotoluene (TFT)		0.106	mg/L	1	0.100	106	84.5 - 115
4-Bromofluorobenzene (4-BFB)		0.0857	$\mathrm{mg/L}$	1	0.100	86	66.5 - 103

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Report Date: September 11, 2009 Work Order: 9082815 Page Number: 19 of 79

HELSTF Chromate Spill Groundwater

Analysis: QC Batch: Prep Batch:	V, Total 63221 53951		Date	lytical Method e Analyzed: ple Preparatio	200	010B 9-09-03 9-09-03		Prep Metho Analyzed B Prepared B	y: RR
		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Vanadi	um	0.0160	0.0160	< 0.000426	mg/L	1	0.000426	0.005	0.000426

Sample: 208427 - HLSF-0143-HMW-037-0809

Laboratory: Lubbock

Analysis: Zn, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951Sample Preparation: 2009-09-03 Prepared By: KV

SDL MQLMethodMQLMDL Based Based Blank Flag Result Result Dilution SDLParameter Result Units (Unadjusted) (Unadjusted) Total Zinc < 0.000465 < 0.00500 < 0.000465 mg/L0.0004650.0050.000465

Method Blank (1)

QC Batch: 63015 Date Analyzed: 2009-08-27 Analyzed By: MD Prep Batch: 53776 QC Preparation: 2009-08-27 Prepared By: MD

Method Blank (1)

4-Bromofluorobenzene (4-BFB)

QC Batch: 63025 Date Analyzed: 2009-08-28 Analyzed By: ER
Prep Batch: 53796 QC Preparation: 2009-08-28 Prepared By: ER

							Reporting
Parameter	Flag		Result		Units		Limits
GRO			< 0.0702		$\mathrm{mg/L}$		0.0702
					G :1	ъ.	
					Spike	$\operatorname{Percent}$	$\operatorname{Recovery}$
$\operatorname{Surrogate}$	Flag	Result	Units	$\operatorname{Dilution}$	${ m Amount}$	Recovery	Limits
Trifluorotoluene (TFT)		0.0979	$_{ m mg/L}$	1	0.100	98	84.5 - 115

mg/L

1

0.100

80

66.5 - 103

0.0803

Report Date: September 11, 2009 Work Order: 9082815 Page Number: 20 of 79

HELSTF Chromate Spill Groundwater

			HELSTF Chrom	ate Spill Gro	oundwater		
Method Blan	nk (1)						
	63042 53812		Date Analyzed: QC Preparation	2009-08-28 a: 2009-08-28			lyzed By: eared By:
Parameter		Flag	Result	-	Units		Reporting Limits
DRO			< 0.876		mg/L		0.876
Surrogate	Flag	Result	Units D	ilution	Spike Amount	Percent Recovery	Recovery Limits
n-Triacontane	0	8.44	m mg/L	1	10.0	84	34.4 - 185
•	nk (1) 63072 53821	Flag	Date Analyzed: QC Preparation:	2009-08-31 2009-08-31 esult	$_{ m Units}$	Analyze Prepare	
Total Mercury	,	Tiag	<0.000		mg/L		3.29e-05
•	nk (1) 63080 53840		Date Analyzed: QC Preparation:	2009-08-31 2009-08-31		Analyze Prepare	d By: AH
Parameter		Flag	Res		Units		$\begin{array}{c} { m Reporting} \\ { m Limits} \end{array}$
Total Cyanide			< 0.01	110	m mg/L		0.011
•	nk (1) 63111 53864		Date Analyzed: QC Preparation:	2009-08-31 2009-08-31		Analyze Prepare	-
Parameter		Flag	Resu	lt	${ m Units}$		$\begin{array}{c} {\rm Reporting} \\ {\rm Limits} \end{array}$

Method Blank (1)

Ammonia-N

QC Batch:	63142	Date Analyzed:	2009-09-01	Analyzed By:	DS
Prep Batch:	53897	QC Preparation:	2009-09-01	Prepared By:	DS

< 0.353

mg/L

0.353

${\it Work~Order:~9082815} \\ {\it HELSTF~Chromate~Spill~Groundwater} \\$

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Parameter	Flag	Result	${ m Units}$	$rac{ m Reporting}{ m Limits}$
HMX	1100	< 0.123	$\mu \mathrm{g/L}$	0.123
RDX		< 0.298	$\mu { m g}/{ m L}$	0.298
1,3,5-Trinitrobenzene		< 0.339	$\mu { m g}/{ m L}$	0.339
1,3-Dinitrobenzene		< 0.389	$\mu { m g}/{ m L}$	0.389
Nitrobenzene		< 0.379	$\mu { m g}/{ m L}$	0.379
Tetryl		< 0.413	$\mu { m g}/{ m L}$	0.413
TNT		< 0.464	$\mu { m g}/{ m L}$	0.464
4-Amino-DNT		< 0.319	$\mu { m g}/{ m L}$	0.319
2-Amino-DNT		< 0.391	$\mu { m g}/{ m L}$	0.391
2,6-DNT		< 0.323	$\mu { m g}/{ m L}$	0.323
2,4-DNT		< 0.366	$\mu { m g}/{ m L}$	0.366
2-NT		< 0.379	$\mu { m g}/{ m L}$	0.379
4-NT		< 0.398	$\mu \mathrm{g}/\mathrm{L}$	0.398
3-NT		< 0.346	$\mu { m g}/{ m L}$	0.346

					Spike	Percent	Recovery
Surrogate	Flag	Result	Units	$\operatorname{Dilution}$	${ m Amount}$	Recovery	Limits
1,2-Dinitrobenzene		2.44	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	98	19.8 - 160

Method Blank (1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

				$\operatorname{Reporting}$
Parameter	Flag	Result	Units	Limits
Total Silver		< 0.00111	$_{ m mg/L}$	0.00111

Method Blank (1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

				$\operatorname{Reporting}$
Parameter	Flag	Result	Units	Limits
Total Aluminum		< 0.00301	m mg/L	0.00301

Method Blank (1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

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Teopore Bacc. Septemoer	,	HELSTF Chromate Spill Ground	water	0
Parameter	Flag	Result	Units	Reporting Limits
Total Arsenic		< 0.00448	mg/L	0.00448
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	Units	Reporting Limits
Total Barium		< 0.00105	m mg/L	0.00105
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	${ m Units}$	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Total Beryllium		< 0.000450	$\mathrm{mg/L}$	0.00045
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Total Cadmium	0	< 0.000303	m mg/L	0.000303
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	${ m Units}$	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Total Cobalt		< 0.000822	m mg/L	0.000822

Method Blank (1)

QC Batch:	63221	Date Analyzed:	2009-09-03	Analyzed By:	RR
Prep Batch:	53951	QC Preparation:	2009-09-03	Prepared By:	KV

Prep Batch: 53951

Work Order: 9082815 HELSTF Chromate Spill Groundwater

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Prepared By: KV

Parameter	Floor	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Total Chromium	Flag	<0.000583	mg/L	0.000583
Total Chromium		₹0.000383	mg/L	0.000983
Method Blank (1)				
QC Batch: 63221		Date Analyzed: 2009-09-03		Analyzed By: RR
Prep Batch: 53951		QC Preparation: 2009-09-03		Prepared By: KV
Parameter	Flag	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Total Copper	riag	<0.000843	mg/L	0.000843
от от		X010000219		3,000013
Method Blank (1)				
QC Batch: 63221		Date Analyzed: 2009-09-03		Analyzed By: RR
Prep Batch: 53951		QC Preparation: 2009-09-03		Prepared By: KV
Parameter	Flag	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Total Iron	Tag	<0.000872	$\frac{\rm mg/L}$	0.000872
Method Blank (1) QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Total Manganese	Flag	<0.000305	$\frac{\rm mg/L}$	0.000305
Method Blank (1)			Oi.	
QC Batch: 63221		Date Analyzed: 2009-09-03		Analyzed By: RR
Prep Batch: 53951		QC Preparation: 2009-09-03		Prepared By: KV
Parameter	Flag	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Total Molybdenum	<u> </u>	< 0.00119	$_{ m mg/L}$	0.00119
Method Blank (1)				
QC Batch: 63221		Date Analyzed: 2009-09-03		Analyzed By: RR

QC Preparation: 2009-09-03

Prep Batch: 53951

Work Order: 9082815 HELSTF Chromate Spill Groundwater

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.11 ~	•		

Prepared By: KV

Ela m	Dogul	Units	Reportin
Flag	Result < 0.00121		Limits 0.00121
	₹0.00121	mg/ L	0.00121
	Date Analyzed: 2009-09-03		Analyzed By: RR
	QC Preparation: 2009-09-03		Prepared By: KV
Ela m	Dogula	IInita	Reportin
Flag			Limits 0.00289
	X0.00203	mg/L	0.00209
	Date Analyzed: 2009-09-03		Analyzed By: RR
	QC Preparation: 2009-09-03		Prepared By: KV
			Reportin
Flag	Rosult	Unite	——————————————————————————————————————
Flag	Result <0.00326	$\begin{array}{c} \text{Units} \\ \text{mg/L} \end{array}$	Limits
	<0.00326 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03	mg/L	Limits 0.00326 Analyzed By: RR Prepared By: KV Reportin
Flag Flag	<pre></pre>	${ m mg/L}$ Units	Limits 0.00326 Analyzed By: RR Prepared By: KV Reportin Limits
	<0.00326 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03	mg/L	Limits 0.00326 Analyzed By: RR Prepared By: KV Reportin
	<pre></pre>	${ m mg/L}$ Units	Analyzed By: RR Prepared By: KV Reportin Limits 0.0044
	<0.00326 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Result <0.00440	${ m mg/L}$ Units	Analyzed By: RR Prepared By: KV Reportin Limits 0.0044 Analyzed By: RR
	<pre></pre>	${ m mg/L}$ Units	Analyzed By: RR Prepared By: KV Reportin Limits 0.0044 Analyzed By: RR
	Flag	Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Flag Result <0.00289 Date Analyzed: 2009-09-03	Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Flag Result Units <0.00289 mg/L Date Analyzed: 2009-09-03

QC Preparation: 2009-09-03

QC Batch:

Prep Batch: 53984

63249

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Analyzed By: SS

Prepared By: SS

D	Tal.	D. L	TT ',	Reporting
Parameter Total Thallium	Flag	Result	Units	Limits
Iotai Inailium		< 0.00488	m mg/L	0.00488
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	Units	Reporting Limits
Total Vanadium		< 0.000426	m mg/L	0.000426
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Total Zinc	Ö	< 0.000465	m mg/L	0.000465
Method Blank (1)				
QC Batch: 63249 Prep Batch: 53984		Date Analyzed: 2009-08-31 QC Preparation: 2009-08-31		Analyzed By: SS Prepared By: SS
Parameter	Flag	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Bromide	1145	<0.192	mg/L	0.192
Method Blank (1)				
QC Batch: 63249		Date Analyzed: 2009-08-31 QC Preparation: 2009-08-31		Analyzed By: SS Prepared By: SS
Prep Batch: 53984				D
Prep Batch: 53984 Parameter Chloride	Flag	Result <0.157	$\frac{\rm Units}{\rm mg/L}$	Reporting Limits 0.157

Date Analyzed:

QC Preparation: 2009-08-31

2009 - 08 - 31

${\it Work~Order:~9082815} \\ {\it HELSTF~Chromate~Spill~Groundwater} \\$

Parameter	Flag	Result	${ m Units}$	$\begin{array}{c} \text{Reporting} \\ \text{Limits} \end{array}$
Fluoride		< 0.204	$\mathrm{mg/L}$	0.204

Method Blank (1)

QC Batch: 63249 Date Analyzed: 2009-08-31 Analyzed By: SS Prep Batch: 53984 QC Preparation: 2009-08-31 Prepared By: SS

				Reporting
Parameter	Flag	Result	${ m Units}$	Limits
Sulfate		< 0.153	m mg/L	0.153

Method Blank (1)

QC Batch: 63261 Date Analyzed: 2009-09-04 Analyzed By: MN Prep Batch: 53996 QC Preparation: 2009-09-02 Prepared By: MN

				Reporting
Parameter	Flag	Result	Units	Limits
Pyridine		< 0.000608	m mg/L	0.000608
N-Nitrosodimethylamine		< 0.000552	m mg/L	0.000552
2-Picoline		< 0.000408	m mg/L	0.000408
Methyl methanesulfonate		< 0.000350	m mg/L	0.00035
Ethyl methanesulfonate		< 0.000448	m mg/L	0.000448
Phenol		< 0.000509	m mg/L	0.000509
Aniline		< 0.000691	m mg/L	0.000691
${ m bis}(2{ m -chloroethyl}){ m ether}$		< 0.000440	m mg/L	0.00044
2-Chlorophenol		< 0.000537	m mg/L	0.000537
1,3-Dichlorobenzene (meta)		< 0.000441	m mg/L	0.000441
1,4-Dichlorobenzene (para)		< 0.000440	m mg/L	0.00044
Benzyl alcohol		< 0.000538	m mg/L	0.000538
1,2-Dichlorobenzene (ortho)		< 0.000443	m mg/L	0.000443
2-Methylphenol		< 0.000726	m mg/L	0.000726
${ m bis}(2{ m -chloroisopropyl}){ m ether}$		< 0.000503	m mg/L	0.000503
4-Methylphenol / 3-Methylphenol		< 0.000512	m mg/L	0.000512
N-Nitrosodi-n-propylamine		< 0.000732	m mg/L	0.000732
$\operatorname{Hexachloroethane}$		< 0.000507	m mg/L	0.000507
$egin{array}{c} { m Acetophenone} \end{array}$		< 0.000424	m mg/L	0.000424
Nitrobenzene		< 0.000465	m mg/L	0.000465
N-Nitrosopiperidine		< 0.000443	$\mathrm{mg/L}$	0.000443
Isophorone		< 0.000619	m mg/L	0.000619
2-Nitrophenol		< 0.000406	m mg/L	0.000406
2,4-Dimethylphenol		< 0.000477	m mg/L	0.000477
${ m bis}(2{ ext{-chloroet}}{ m hoxy}){ m methane}$		< 0.000432	m mg/L	0.000432
2,4-Dichlorophenol		< 0.000400	m mg/L	0.0004
1,2,4-Trichlorobenzene		< 0.000404	$\mathrm{mg/L}$	0.000404

 $continued \dots$

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method blank continued . . .

meenou outen communeu				Reporting
Parameter	Flag	Result	Units	Limits
Benzoic acid		< 0.00163	m mg/L	0.00163
Naphthalene		< 0.000489	m mg/L	0.000489
${ m a,a-Dimethyl}$ phenethylamine		< 0.00129	$\mathrm{mg/L}$	0.00129
4-Chloroaniline		< 0.000378	$\mathrm{mg/L}$	0.000378
2,6-Dichlorophenol		< 0.000484	${ m mg/L}$	0.000484
${\it Hexachlorobut}$ adiene		< 0.000517	$\mathrm{mg/L}$	0.000517
N-Nitroso-di-n-butylamine		< 0.000656	$\mathrm{mg/L}$	0.000656
4-Chloro-3-methylphenol		< 0.000522	$\mathrm{mg/L}$	0.000522
$2 ext{-Methylnaphthalene}$		< 0.000423	$\mathrm{mg/L}$	0.000423
$1 ext{-} ext{Methylnapht}$ halene		< 0.000495	$\mathrm{mg/L}$	0.000495
1,2,4,5-Tetrachlorobenzene		< 0.000612	$\mathrm{mg/L}$	0.000612
${\it Hexachlorocyclopenta}$ diene		< 0.000558	$\mathrm{mg/L}$	0.000558
2,4,6-Trichlorophenol		< 0.000794	${ m mg/L}$	0.000794
2,4,5-Trichlorophenol		< 0.000834	${ m mg/L}$	0.000834
$2 ext{-Chloronaphthalene}$		< 0.000416	$\mathrm{mg/L}$	0.000416
1-Chloronaphthalene		< 0.000476	$\mathrm{mg/L}$	0.000476
2-Nitroaniline		< 0.000760	$\mathrm{mg/L}$	0.00076
${\bf Dimethylphthalate}$		< 0.000643	$\mathrm{mg/L}$	0.000643
${ m Acenapht hylene}$		< 0.000586	$\mathrm{mg/L}$	0.000586
2,6-Dinitrotoluene		< 0.000640	$\mathrm{mg/L}$	0.00064
3-Nitroaniline		< 0.000721	$\mathrm{mg/L}$	0.000721
${ m Acenaphthene}$		< 0.000423	${ m mg/L}$	0.000423
2,4-Dinitrophenol		< 0.000220	${ m mg/L}$	0.00022
Dibenzofuran		< 0.000408	m mg/L	0.000408
Pentachlorobenzene		< 0.000571	m mg/L	0.000571
4-Nitrophenol		< 0.00185	m mg/L	0.00185
2,4-Dinitrotoluene		< 0.000911	m mg/L	0.000911
$1 ext{-Naphthylamine}$		< 0.000688	m mg/L	0.000688
2,3,4,6-Tetrachlorophenol		< 0.000565	m mg/L	0.000565
2-Naphthylamine		< 0.000699	${ m mg/L}$	0.000699
Fluorene		< 0.000648	m mg/L	0.000648
4-Chlorophenyl-phenylether		< 0.000619	m mg/L	0.000619
${ m Diethylphthalate}$		< 0.000828	${ m mg/L}$	0.000828
4-Nitroaniline		< 0.000702	${ m mg/L}$	0.000702
Diphenylhydrazine		< 0.000657	${ m mg/L}$	0.000657
4,6-Dinitro- 2 -methylphenol		< 0.00198	m mg/L	0.00198
Diphenylamine		< 0.000440	m mg/L	0.00044
4-Bromophenyl-phenylether		< 0.000550	${ m mg/L}$	0.00055
Phenacetin		< 0.000605	${ m mg/L}$	0.000605
Hexachlorobenzene		< 0.000506	${ m mg/L}$	0.000506
4-Aminobiphenyl		< 0.000527	${ m mg/L}$	0.000527
Pentachlorophenol		< 0.000435	m mg/L	0.000435
${ m Anthracene}$		< 0.000428	m mg/L	0.000428
${\bf Pentach loron it robenzene}$		< 0.000408	m mg/L	0.000408
Pronamide		< 0.000476	m mg/L	0.000476
Phenanthrene		< 0.000548	m mg/L	0.000548
Di-n-butylphthalate		< 0.000483	m mg/L	0.000483

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method	hlank	continued

				Reporting
Parameter	Flag	Result	Units	Limits
Fluoranthene		< 0.000632	m mg/L	0.000632
Benzidine		< 0.00238	m mg/L	0.00238
Pyrene		< 0.000723	m mg/L	0.000723
p-Dimethylaminoazobenzene		< 0.000902	m mg/L	0.000902
Butylbenzylphthalate		< 0.000445	m mg/L	0.000445
$\operatorname{Benzo}(\operatorname{a})\operatorname{anthracene}$		< 0.000527	m mg/L	0.000527
3,3-Dichlorobenzidine		< 0.00118	m mg/L	0.00118
Chrysene		< 0.000638	m mg/L	0.000638
bis(2-ethylhexyl)phthalate		< 0.000561	m mg/L	0.000561
Di-n-octylphthalate		< 0.00116	m mg/L	0.00116
Benzo(b)fluoranthene		< 0.000879	m mg/L	0.000879
Benzo(k)fluoranthene		< 0.000845	m mg/L	0.000845
7,12-Dimethylbenz (a) anthracene		< 0.00102	m mg/L	0.00102
Benzo(a)pyrene		< 0.00167	m mg/L	0.00167
3-Methylcholanthrene		< 0.000908	m mg/L	0.000908
${ m Dibenzo(a,j)}$ acridine		< 0.00129	m mg/L	0.00129
Indeno(1,2,3-cd)pyrene		< 0.000862	m mg/L	0.000862
${ m Dibenzo}({ m a,h}) { m anthracene}$		< 0.000809	m mg/L	0.000809
$\operatorname{Benzo}(g,h,i)$ perylene		< 0.000949	m mg/L	0.000949

					Spike	$\operatorname{Percent}$	$\operatorname{Recovery}$
Surrogate	Flag	Result	Units	Dilution	${f Amount}$	Recovery	Limits
2-Fluorophenol		0.0250	$\mathrm{mg/L}$	1	0.0800	31	10 - 53.1
Phenol-d5		0.0158	$\mathrm{mg/L}$	1	0.0800	20	10 - 36.9
${ m Nitrobenzene-d5}$		0.0378	$\mathrm{mg/L}$	1	0.0800	47	23.8 - 108
2-Fluorobiphenyl		0.0396	$_{ m mg/L}$	1	0.0800	50	15.9 - 127
2,4,6-Tribromophenol		0.0493	$\mathrm{mg/L}$	1	0.0800	62	10 - 123
Terphenyl-d14		0.0437	m mg/L	1	0.0800	55	17.2 - 160

Method Blank (1)

QC Batch: 63272 Date Analyzed: 2009-09-01 Analyzed By: MD Prep Batch: 54007 QC Preparation: 2009-09-01 Prepared By: MD

				$\operatorname{Reporting}$
Parameter	Flag	Result	Units	Limits
Total Dissolved Solids		< 5.00	$\mathrm{mg/L}$	5

Method Blank (1)

QC Batch: 63291 Date Analyzed: 2009-09-02 Analyzed By: AH Prep Batch: 54023 QC Preparation: 2009-09-02 Prepared By: AH

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Parameter	F	lag	Result	Units	-	$rac{ ext{orting}}{ ext{mits}}$
Total Kjeldahl Nitrogen - N		0	< 2.45	m mg/L		.45
Method Blank (1)						
QC Batch: 63340		Date Analyzed:	2009-09-08		Analyzed By:	KV
Prep Batch: 54065	(QC Preparation:	2009-09-08		Prepared By:	KV
D	T 1	_	S 1	TT 1:	•	orting
Parameter Total Organic Carbon	Flag	1	Result 0.848	$\frac{\rm Units}{\rm mg/L}$		$\frac{\text{mits}}{401}$
Total Olganic Carbon			0.010	6/12	0.	101
Method Blank (1)						
QC Batch: 63355		Date Analyzed:	2009-09-09		Analyzed By:	RR
Prep Batch: 53951	(QC Preparation:	2009-09-03		Prepared By:	KV
Parameter	Flag	Resi	11+	Units		$_{ m orting}$
Total Calcium	riag	<0.1		mg/L		$\frac{117}{117}$
Method Blank (1) QC Batch: 63355 Prep Batch: 53951		Oate Analyzed: QC Preparation:	2009-09-09 2009-09-03		Analyzed By: Prepared By:	RR KV
Parameter	Flag	Res	l+	Units	-	$rac{ m orting}{ m mits}$
Total Potassium	Tag	<0.		mg/L		$\frac{1172}{172}$
Method Blank (1)						
QC Batch: 63355		Date Analyzed:	2009-09-09		Analyzed By:	RR
Prep Batch: 53951	(QC Preparation:	2009-09-03		Prepared By:	KV
Parameter	Flag	Ro	sult	Units		$rac{ ext{orting}}{ ext{mits}}$
Total Magnesium	Flag		160	$\frac{\rm mg/L}$.16
Method Blank (1)						
QC Batch: 63355		Date Analyzed:	2009-09-09		Analyzed By:	RR
Prep Batch: 53951	(QC Preparation:	2009-09-03		Prepared By:	KV

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Parameter	Flag]+	${ m Units}$		$ \begin{array}{c} \text{orting} \\ \text{mits} \end{array} $	
Total Sodium	rag	Res < 0.05		$\frac{\rm mg/L}$.05
10vai poqidin		₹0.0€		mg/ L	0	.00
Method Blank (1)						
QC Batch: 63379		Date Analyzed:	2009-09-08		Analyzed By:	$_{ m JG}$
Prep Batch: 54095		QC Preparation:	2009-09-08		Prepared By:	$_{ m JG}$
.	T)	ъ.			_	orting
Parameter Hydroxide Alkalinity	Flag	Resul		$\frac{\text{Units}}{\text{mg/L as CaCo3}}$		mits
Carbonate Alkalinity		<1.0 <1.0		mg/L as CaCo3		1 1
Bicarbonate Alkalinity		<4.0		mg/L as CaCo3		4
Total Alkalinity		<4.0		mg/L as CaCo3		$\overline{4}$
Method Blank (1)						
QC Batch: 63411 Prep Batch: 54129		Date Analyzed: QC Preparation:	2009-09-09 2009-09-08		Analyzed By: Prepared By:	MD MD
					D	, •
Parameter	Flag	Res	ılt	Units	_	$rac{1}{2} \operatorname{corting} $
Oil and Grease	1145	<6.		$\frac{\rm mg/L}$		3.6
				Oi .		
Method Blank (1)						
QC Batch: 63461		Date Analyzed:	2009-09-11		Analyzed By:	RR
Prep Batch: 54153		QC Preparation:	2009-09-11		Prepared By:	KV
					Rep	orting
Parameter	Flag		Result	Units	Lii	$_{ m mits}$
Parameter Dissolved Chromium	Flag	<0	Result .000583	$\begin{array}{c} \rm Units \\ \rm mg/L \end{array}$	Lii	_
	Flag	<0			Lii	$_{ m mits}$
Dissolved Chromium Method Blank (1)	Flag		.000583		Li ₁	mits 00583
Dissolved Chromium	Flag	<0 Date Analyzed: QC Preparation:			Lii	$_{ m mits}$
Dissolved Chromium Method Blank (1) QC Batch: 63485		Date Analyzed: QC Preparation:	2009-09-11 2009-09-11		Analyzed By: Prepared By: Rep	mits 00583 KV KV KV orting
Dissolved Chromium Method Blank (1) QC Batch: 63485	Flag Fl	Date Analyzed: QC Preparation: ag	2009-09-11		Analyzed By: Prepared By: Rep	mits 00583 KV KV

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Duplicate (1) Duplicated San

QC Batch: 63082 Date Analyzed: 2009-08-27 Prep Batch: 53841 QC Preparation: 2009-08-27

Analyzed By: JG Prepared By: $_{
m JG}$

	Duplicate	\mathbf{Sample}				RPD
Param	Result	Result	Units	Dilution	RPD	Limit
pН	7.33	7.33	s.u.	1	0	1.1

Duplicate (1) Duplicated Sample: 208087

QC Batch: 63272 Date Analyzed: 2009-09-01 Analyzed By: MD Prep Batch: 54007 QC Preparation: 2009-09-01 Prepared By: MD

	Duplicate	\mathbf{Sample}				RPD
Param	Result	Result	Units	Dilution	RPD	Limit
Total Dissolved Solids	16000	15800	$_{ m mg/L}$	1	1	10

Duplicate (1) Duplicated Sample: 208085

QC Batch: Date Analyzed: 2009-09-08 63379 Analyzed By: JG Prep Batch: 54095 QC Preparation: 2009-09-08 Prepared By: JG

Param	Duplicate Result	$\begin{array}{c} { m Sample} \\ { m Result} \end{array}$	Units	Dilution	RPD	$egin{array}{c} ext{RPD} \ ext{Limit} \end{array}$
raram	nesuit	\mathbf{nesum}	Umus	Dilution	$\mathbf{n}_{\mathbf{L}}\mathbf{D}$	PIIIII
Hydroxide Alkalinity	< 1.00	<1.00	mg/L as CaCo3	1	0	20
Carbonate Alkalinity	< 1.00	< 1.00	mg/L as $CaCo3$	1	0	20
Bicarbonate Alkalinity	164	160	mg/L as $CaCo3$	1	2	20
Total Alkalinity	164	160	mg/L as CaCo3	1	2	20

Laboratory Control Spike (LCS-1)

Date Analyzed: QC Batch: 630252009-08-28 Analyzed By: ER Prep Batch: 53796 QC Preparation: 2009-08-28 Prepared By: ER

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit
$\overline{\mathrm{GRO}}$	0.954	mg/L	1	1.00	< 0.0702	95	74.6 - 119

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	$\mathrm{Rec}.$	Limit	RPD	Limit
$\overline{\mathrm{GRO}}$	1.02	mg/L	1	1.00	< 0.0702	102	74.6 - 119	6	20

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	LCS	LCSD			Spike	LCS	LCSD	Rec .
Surrogate	Result	Result	Units	Dil.	${ m Amount}$	Rec .	Rec .	Limit
Trifluorotoluene (TFT)	0.0993	0.0977	mg/L	1	0.100	99	98	77.8 - 114
4-Bromofluorobenzene (4-BFB)	0.0888	0.0853	$_{ m mg/L}$	1	0.100	89	85	80.3 - 111

Laboratory Control Spike (LCS-1)

QC Batch: 63042 Date Analyzed: 2009-08-28 Analyzed By: Prep Batch: 53812 QC Preparation: 2009-08-28 Prepared By:

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
DRO	31.0	$\mathrm{mg/L}$	1	25.0	< 0.876	124	74.3 - 158

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit	RPD	Limit
DRO	30.0	mg/L	1	25.0	< 0.876	120	74.3 - 158	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCS	LCSD			Spike	LCS	LCSD	$\mathrm{Rec}.$
Surrogate	Result	Result	Units	Dil.	${f Amount}$	Rec.	Rec .	Limit
n-Triacont ane	8.97	8.78	mg/L	1	10.0	90	88	34.4 - 149

Laboratory Control Spike (LCS-1)

QC Batch: 63072 Date Analyzed: 2009-08-31 Analyzed By: TP
Prep Batch: 53821 QC Preparation: 2009-08-31 Prepared By: TP

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec}.$	Limit
Total Mercury	0.000930	$\mathrm{mg/L}$	1	0.00100	< 0.0000329	93	90.3 - 108

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		Rec .		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit	RPD	Limit
Total Mercury	0.000940	mg/L	1	0.00100	< 0.0000329	94	90.3 - 108	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63142 Date Analyzed: 2009-09-01 Analyzed By: DS
Prep Batch: 53897 QC Preparation: 2009-09-01 Prepared By: DS

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	LCS			Spike	Matrix		Rec.
Param	Result	Units	Dil.	${ m Amount}$	Result	$\mathrm{Rec.}$	Limit
HMX	2.49	$\mu \mathrm{g/L}$	1	2.50	< 0.123	100	63.5 - 125
RDX	2.34	$\mu { m g}/{ m L}$	1	2.50	< 0.298	94	74.5 - 124
1,3,5-Trinitrobenzene	2.37	$\mu { m g}/{ m L}$	1	2.50	< 0.339	95	54.1 - 131
1,3-Dinitrobenzene	2.45	$\mu { m g}/{ m L}$	1	2.50	< 0.389	98	72 - 112
Nitrobenzene	2.49	$\mu { m g}/{ m L}$	1	2.50	< 0.379	100	72.5 - 126
Tetryl	2.27	$\mu { m g}/{ m L}$	1	2.50	< 0.413	91	35.9 - 149
TNT	2.45	$\mu { m g}/{ m L}$	1	2.50	< 0.464	98	40.7 - 129
4-Amino-DNT	2.41	$\mu { m g}/{ m L}$	1	2.50	< 0.319	96	80 - 120
2-Amino-DNT	2.50	$\mu { m g}/{ m L}$	1	2.50	< 0.391	100	80 - 120
2,6-DNT	2.05	$\mu { m g}/{ m L}$	1	2.50	< 0.323	82	80 - 120
2,4-DNT	2.48	$\mu { m g}/{ m L}$	1	2.50	< 0.366	99	80 - 120
2-NT	2.33	$\mu { m g}/{ m L}$	1	2.50	< 0.379	93	49.8 - 139
4-NT	2.26	$\mu { m g}/{ m L}$	1	2.50	< 0.398	90	56.3 - 141
3-NT	2.19	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.346	88	66.2 - 129

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	$\mathrm{Rec}.$	Limit	RPD	Limit
HMX	2.58	$\mu \mathrm{g/L}$	1	2.50	< 0.123	103	63.5 - 125	4	20
RDX	2.49	$\mu { m g/L}$	1	2.50	< 0.298	100	74.5 - 124	6	20
1,3,5-Trinitrobenzene	2.40	$\mu { m g/L}$	1	2.50	< 0.339	96	54.1 - 131	1	20
1,3-Dinitrobenzene	2.47	$\mu { m g/L}$	1	2.50	< 0.389	99	72 - 112	1	20
Nitrobenzene	2.49	$\mu { m g/L}$	1	2.50	< 0.379	100	72.5 - 126	0	20
Tetryl	2.35	$\mu { m g/L}$	1	2.50	< 0.413	94	35.9 - 149	4	20
TNT	2.52	$\mu { m g/L}$	1	2.50	< 0.464	101	40.7 - 129	3	20
4-Amino-DNT	2.51	$\mu { m g/L}$	1	2.50	< 0.319	100	80 - 120	4	20
2-Amino-DNT	2.70	$\mu { m g/L}$	1	2.50	< 0.391	108	80 - 120	8	20
2,6-DNT	2.31	$\mu \mathrm{g/L}$	1	2.50	< 0.323	92	80 - 120	12	20
2,4-DNT	2.67	$\mu { m g}/{ m L}$	1	2.50	< 0.366	107	80 - 120	7	20
2-NT	2.37	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.379	95	49.8 - 139	2	20
4-NT	2.15	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.398	86	56.3 - 141	5	20
3-NT	2.45	$\mu { m g}/{ m L}$	1	2.50	< 0.346	98	66.2 - 129	11	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCS	LCSD			Spike	LCS	LCSD	${ m Rec.}$
Surrogate	Result	Result	Units	Dil.	${f Amount}$	Rec.	$\mathrm{Rec}.$	Limit
1,2-Dinitrobenzene	2.39	2.34	$\mu \mathrm{g/L}$	1	2.50	96	94	53 - 134

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit
Total Silver	0.124	${ m mg/L}$	1	0.125	< 0.00111	99	85 - 115

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit	RPD	Limit
Total Silver	0.124	mg/L	1	0.125	< 0.00111	99	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil .	${f Amount}$	Result	Rec .	Limit
Total Aluminum	0.953	$\mathrm{mg/L}$	1	1.00	< 0.00301	95	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Aluminum	0.951	mg/L	1	1.00	< 0.00301	95	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Arsenic	0.478	mg/L	1	0.500	< 0.00448	96	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	$\mathbf{A}\mathbf{mount}$	Result	Rec.	Limit	RPD	Limit
Total Arsenic	0.476	mg/L	1	0.500	< 0.00448	95	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03

Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Total Barium	1.03	${ m mg/L}$	1	1.00	< 0.00105	103	85 - 115

	T CCD			G 11	3.5		T.		DDD
	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	Amount	Result	Rec .	Limit	RPD	Limit
Total Barium	1.03	mg/L	1	1.00	< 0.00105	103	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Beryllium	0.0248	$\mathrm{mg/L}$	1	0.0250	< 0.000450	99	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	Amount	Result	Rec .	Limit	RPD	Limit
Total Beryllium	0.0256	$\mathrm{mg/L}$	1	0.0250	< 0.000450	102	85 - 115	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Cadmium	0.250	m mg/L	1	0.250	< 0.000303	100	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	A mount	Result	Rec.	Limit	RPD	Limit
Total Cadmium	0.250	mg/L	1	0.250	< 0.000303	100	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03

Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Cobalt	0.246	m mg/L	1	0.250	< 0.000822	98	85 - 115

	LCSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Cobalt	0.246	mg/L	1	0.250	< 0.000822	98	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

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	$_{ m LCS}$			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Total Chromium	0.0985	$\mathrm{mg/L}$	1	0.100	< 0.000583	98	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Total Chromium	0.0980	$\mathrm{mg/L}$	1	0.100	< 0.000583	98	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		Rec.
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Copper	0.131	mg/L	1	0.125	< 0.000843	105	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil .	$\mathbf{A}\mathbf{mount}$	Result	Rec.	Limit	RPD	Limit
Total Copper	0.133	mg/L	1	0.125	< 0.000843	106	85 - 115	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Iron	0.570	$\mathrm{mg/L}$	1	0.500	< 0.000872	114	85 - 115

	LCSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Total Iron	0.516	mg/L	1	0.500	< 0.000872	103	85 - 115	10	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${ m Amount}$	Result	$\mathrm{Rec}.$	Limit
Total Manganese	0.253	$\mathrm{mg/L}$	1	0.250	< 0.000305	101	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec.}$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Manganese	0.254	mg/L	1	0.250	< 0.000305	102	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil .	${f Amount}$	Result	Rec.	Limit
Total Molvbdenum	0.544	mg/L	1	0.500	< 0.00119	109	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Molybdenum	0.543	mg/L	1	0.500	< 0.00119	109	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

 QC Batch:
 63221
 Date Analyzed:
 2009-09-03

 Prep Batch:
 53951
 QC Preparation:
 2009-09-03

Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Total Nickel	0.252	${ m mg/L}$	1	0.250	< 0.00121	101	85 - 115

	T CCD			G 11	3.5		ъ		DDD
	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	A mount	Result	$\mathrm{Rec}.$	Limit	RPD	Limit
Total Nickel	0.251	mg/L	1	0.250	< 0.00121	100	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Phosphorous	0.473	$\mathrm{mg/L}$	1	0.500	< 0.00289	95	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec.}$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Phosphorous	0.470	$\mathrm{mg/L}$	1	0.500	< 0.00289	94	85 - 115	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Lead	0.468	$\mathrm{mg/L}$	1	0.500	< 0.00326	94	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit	RPD	Limit
Total Lead	0.464	mg/L	1	0.500	< 0.00326	93	85 - 115	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Antimony	0.249	$\mathrm{mg/L}$	1	0.250	< 0.00440	100	85 - 115

	LCSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Total Antimony	0.248	mg/L	1	0.250	< 0.00440	99	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	$\mathrm{Dil}.$	${f Amount}$	Result	Rec .	Limit
Total Selenium	0.442	$\mathrm{mg/L}$	1	0.500	< 0.00508	88	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Selenium	0.443	mg/L	1	0.500	< 0.00508	89	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Thallium	0.503	$\mathrm{mg/L}$	1	0.500	< 0.00488	101	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Thallium	0.503	mg/L	1	0.500	< 0.00488	101	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Vanadium	0.246	$\mathrm{mg/L}$	1	0.250	< 0.000426	98	85 - 115

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	A mount	Result	Rec.	Limit	RPD	Limit
Total Vanadium	0.248	mg/L	1	0.250	< 0.000426	99	85 - 115	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Total Zinc	0.251	$\mathrm{mg/L}$	1	0.250	< 0.000465	100	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	Amount	Result	Rec .	Limit	RPD	Limit
Total Zinc	0.253	mg/L	1	0.250	< 0.000465	101	85 - 115	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63249 Prep Batch: 53984 Date Analyzed: 2009-08-31 QC Preparation: 2009-08-31 Analyzed By: SS Prepared By: SS

	LCS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Bromide	5.06	$\mathrm{mg/L}$	1	5.00	< 0.192	101	90 - 110

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		Rec.		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Bromide	4.95	mg/L	1	5.00	< 0.192	99	90 - 110	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63249 Prep Batch: 53984 Date Analyzed: 2009-08-31 QC Preparation: 2009-08-31

Analyzed By: SS Prepared By: SS

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Chloride	23.8	m mg/L	1	25.0	< 0.157	95	90 - 110

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SS

	LCSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Chloride	23.9	mø/L	1	25.0	< 0.157	96	90 - 110	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63249 Date Analyzed: 2009-08-31 Analyzed By: Prep Batch: 53984 QC Preparation: 2009-08-31 Prepared By:

LCS Spike Rec. Matrix Param Result Units Dil. Amount Result Rec. Limit Fluoride 5.44 5.00 < 0.204109 90 - 110 mg/L1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil .	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Fluoride	5.38	mg/L	1	5.00	< 0.204	108	90 - 110	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63249 Date Analyzed: 2009-08-31 Analyzed By: SS Prep Batch: 53984 QC Preparation: 2009-08-31 Prepared By: SS

LCS Spike Matrix Rec. Param Result Units Dil. Limit Amount Result Rec. Sulfate 24.2 mg/L25.0< 0.153 97 90 - 110

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		Rec.		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Sulfate	24.1	mg/L	1	25.0	< 0.153	96	90 - 110	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch:63261Date Analyzed:2009-09-04Analyzed By:MNPrep Batch:53996QC Preparation:2009-09-02Prepared By:MN

	LCS			Spike	Matrix		Rec .
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Phenol	0.0205	$\mathrm{mg/L}$	1	0.0800	< 0.000509	26	10 - 66.5
2-Chlorophenol	0.0411	$\mathrm{mg/L}$	1	0.0800	< 0.000537	51	11.2 - 108
1,4-Dichlorobenzene (para)	0.0378	$\mathrm{mg/L}$	1	0.0800	< 0.000440	47	16 - 101

 $continued \dots$

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$control\ spikes\ continued\ \dots$								
		$_{ m LCS}$			Spike	Matrix		Rec .
Param		Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
N-Nitrosodi-n-propylamine		0.0474	mg/L	1	0.0800	< 0.000732	59	10 - 142
1,2,4-Trichlorobenzene		0.0389	$\mathrm{mg/L}$	1	0.0800	< 0.000404	49	18 - 118
Naphthalene		0.0410	$\mathrm{mg/L}$	1	0.0800	< 0.000489	51	20.2 - 114
4-Chloro-3-methylphenol		0.0507	$\mathrm{mg/L}$	1	0.0800	< 0.000522	63	21.5 - 125
${ m Acenapht hylene}$		0.0474	${ m mg/L}$	1	0.0800	< 0.000586	59	25.8 - 121
${ m Acenaphthene}$		0.0474	${ m mg/L}$	1	0.0800	< 0.000423	59	33.5 - 122
4-Nitrophenol		0.0191	${ m mg/L}$	1	0.0800	< 0.00185	24	10 - 125
2,4-Dinitrotoluene		0.0543	${ m mg/L}$	1	0.0800	< 0.000911	68	53 - 130
Fluorene		0.0494	${ m mg/L}$	1	0.0800	< 0.000648	62	44.6 - 117
${ m Pentachlorophenol}$		0.0373	${ m mg/L}$	1	0.0800	< 0.000435	47	10 - 139
${ m Anthracene}$		0.0511	${ m mg/L}$	1	0.0800	< 0.000428	64	57.5 - 115
Phenanthrene		0.0509	${ m mg/L}$	1	0.0800	< 0.000548	64	55.5 - 118
Fluoranthene		0.0545	${ m mg/L}$	1	0.0800	< 0.000632	68	57 - 122
Pyrene		0.0468	${ m mg/L}$	1	0.0800	< 0.000723	58	58.5 - 130
$\mathrm{Benzo}(\mathrm{a}) \mathrm{anthracene}$	1	0.0463	${ m mg/L}$	1	0.0800	< 0.000527	58	63.4 - 109
Chrysene		0.0508	${ m mg/L}$	1	0.0800	< 0.000638	64	54.7 - 114
${ m Benzo(b)}$ fluoranthene	2	0.0477	${ m mg/L}$	1	0.0800	< 0.000879	60	64.8 - 120
$\mathrm{Benzo}(\mathtt{k}) \mathrm{fluoranthene}$		0.0619	$\mathrm{mg/L}$	1	0.0800	< 0.000845	77	70.3 - 114
$\mathrm{Benzo}(\mathrm{a})\mathrm{pyrene}$		0.0575	${ m mg/L}$	1	0.0800	< 0.00167	72	63.7 - 120
Indeno(1,2,3-cd)pyrene		0.0568	${ m mg/L}$	1	0.0800	< 0.000862	71	65.4 - 119
${ m Dibenzo}({ m a,h}) { m anthracene}$		0.0572	${ m mg/L}$	1	0.0800	< 0.000809	72	68.7 - 117
$\operatorname{Benzo}(g,h,i)$ perylene		0.0598	m mg/L	1	0.0800	< 0.000949	75	57.2 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Phenol	0.0210	mg/L	1	0.0800	< 0.000509	26	10 - 66.5	2	20
2-Chlorophenol	0.0432	mg/L	1	0.0800	< 0.000537	54	11.2 - 108	5	20
1,4-Dichlorobenzene (para)	0.0398	mg/L	1	0.0800	< 0.000440	50	16 - 101	5	20
N-Nitrosodi-n-propylamine	0.0506	$\mathrm{mg/L}$	1	0.0800	< 0.000732	63	10 - 142	6	20
1,2,4-Trichlorobenzene	0.0413	$\mathrm{mg/L}$	1	0.0800	< 0.000404	52	18 - 118	6	20
Naphthalene	0.0436	mg/L	1	0.0800	< 0.000489	54	20.2 - 114	6	20
4-Chloro-3-methylphenol	0.0534	mg/L	1	0.0800	< 0.000522	67	21.5 - 125	5	20
${f Acenaphthylene}$	0.0515	mg/L	1	0.0800	< 0.000586	64	25.8 - 121	8	20
${f Acenaphthene}$	0.0505	$\mathrm{mg/L}$	1	0.0800	< 0.000423	63	33.5 - 122	6	20
4-Nitrophenol	0.0220	mg/L	1	0.0800	< 0.00185	28	10 - 125	14	20
2,4-Dinitrotoluene	0.0577	mg/L	1	0.0800	< 0.000911	72	53 - 130	6	20
$\operatorname{Fluorene}$	0.0533	$\mathrm{mg/L}$	1	0.0800	< 0.000648	67	44.6 - 117	8	20
${ m Pentachlorophenol}$	0.0388	$\mathrm{mg/L}$	1	0.0800	< 0.000435	48	10 - 139	4	20
${ m Anthracene}$	0.0530	$\mathrm{mg/L}$	1	0.0800	< 0.000428	66	57.5 - 115	4	20
Phenanthrene	0.0521	$\mathrm{mg/L}$	1	0.0800	< 0.000548	65	55.5 - 118	2	20
Fluoranthene	0.0564	mg/L	1	0.0800	< 0.000632	70	57 - 122	3	20
Pyrene	0.0495	mg/L	1	0.0800	< 0.000723	62	58.5 - 130	6	20
$\operatorname{Benzo}(\operatorname{a}) \operatorname{anthracene}$	0.0486	$\mathrm{mg/L}$	1	0.0800	< 0.000527	61	63.4 - 109	5	20

 $continued \dots$

 $^{^1\}mathrm{Spike}$ analyte out of control limits. Results biased low. • $^2\mathrm{Spike}$ analyte out of control limits. Results biased low. • $^3\mathrm{Spike}$ analyte out of control limits. Results biased low. •

Benzo(g,h,i)perylene

Work Order: 9082815 HELSTF Chromate Spill Groundwater

0.0800

< 0.000949

80

57.2 - 125

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7

20

control spikes continued . . . LCSD Spike Matrix Rec. RPD Param Result Units Dil. Amount Result Limit RPD Rec. Limit Chrysene 0.0541mg/L1 0.0800 < 0.000638 68 54.7 - 114 6 20 Benzo(b)fluoranthene mg/L0.080064.8 - 120 20 0.05331 < 0.000879 67 11 Benzo(k)fluoranthene 0.0696mg/L0.0800< 0.00084587 70.3 - 11412 20 1 Benzo(a)pyrene 63.7 - 1209 20 0.0630mg/L1 0.0800< 0.0016779 Indeno(1,2,3-cd)pyrene 65.4 - 11920 0.0615mg/L0.0800< 0.00086277 8 1 68.7 - 117Dibenzo(a,h)anthracene 0.0630mg/L1 0.0800< 0.000809 79 10 20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

1

mg/L

0.0643

	LCS	LCSD			Spike	LCS	LCSD	Rec .
Surrogate	Result	Result	Units	Dil.	${f Amount}$	Rec .	$\mathrm{Rec}.$	Limit
2-Fluorophenol	0.0260	0.0278	$\mathrm{mg/L}$	1	0.0800	32	35	10 - 53.1
${ m Phenol-d5}$	0.0198	0.0210	${ m mg/L}$	1	0.0800	25	26	10 - 36.9
${ m Nitrobenzene-d5}$	0.0419	0.0439	mg/L	1	0.0800	52	55	23.8 - 108
2-Fluorobiphenyl	0.0429	0.0466	mg/L	1	0.0800	54	58	15.9 - 127
2,4,6-Tribromophenol	0.0565	0.0609	mg/L	1	0.0800	71	76	10 - 123
Terphenyl-d14	0.0461	0.0494	m mg/L	1	0.0800	58	62	17.2 - 160

Laboratory Control Spike (LCS-1)

QC Batch: 63340 Date Analyzed: 2009-09-08 Analyzed By: KV Prep Batch: 54065 QC Preparation: 2009-09-08 Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Total Organic Carbon	48.4	mg/L	1	50.0	< 0.401	97	89.5 - 114

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit	RPD	Limit
Total Organic Carbon	49.4	mg/L	1	50.0	< 0.401	99	89.5 - 114	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

	LCS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${ m Amount}$	Result	$\mathrm{Rec.}$	Limit
Total Calcium	54.2	mg/L	1	50.0	< 0.117	108	85 - 115

	LCSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil .	Amount	Result	Rec.	Limit	RPD	Limit
Total Calcium	50.5	mg/L	1	50.0	< 0.117	101	85 - 115	7	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63355 Prep Batch: 53951 Date Analyzed: 2009-09-09 QC Preparation: 2009-09-03

Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Potassium	53.3	$\mathrm{mg/L}$	1	50.0	< 0.172	107	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil .	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Potassium	50.2	$\mathrm{mg/L}$	1	50.0	< 0.172	100	85 - 115	6	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63355 Prep Batch: 53951 Date Analyzed: 2009-09-09 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${\bf Amount}$	Result	$\mathrm{Rec}.$	Limit
Total Magnesium	52.2	mg/L	1	50.0	< 0.160	104	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		Rec.		RPD
Param	Result	Units	Dil .	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Magnesium	50.1	$\mathrm{mg/L}$	1	50.0	< 0.160	100	85 - 115	4	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63355 Prep Batch: 53951 Date Analyzed: 2009-09-09 QC Preparation: 2009-09-03

Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Total Sodium	53.2	$\mathrm{mg/L}$	1	50.0	< 0.0500	106	85 - 115

Work Order: 9082815 HELSTF Chromate Spill Groundwater

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Sodium	50.6	mg/L	1	50.0	< 0.0500	101	85 - 115	5	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63411 Prep Batch: 54129 Date Analyzed: 2009-09-09 QC Preparation: 2009-09-08

Analyzed By: MD Prepared By: MD

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		LCS			Spike	Matrix		$\mathrm{Rec}.$
Param		Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Oil and Grease	4	18.3	m mg/L	1	40.0	< 3.60	46	78 - 114

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Oil and Grease	5 17.4	$\mathrm{mg/L}$	1	40.0	< 3.60	44	78 - 114	5	18

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63461 Prep Batch: 54153 Date Analyzed: 2009-09-11 QC Preparation: 2009-09-11 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Dissolved Chromium	0.0990	$\mathrm{mg/L}$	1	0.100	< 0.000583	99	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	$_{ m LCSD}$			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	A mount	Result	Rec.	Limit	RPD	Limit
Dissolved Chromium	0.0950	mg/L	1	0.100	< 0.000583	95	85 - 115	4	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208427

 QC Batch:
 63015
 Date Analyzed:
 2009-08-27

 Prep Batch:
 53776
 QC Preparation:
 2009-08-27

Analyzed By: MD Prepared By: MD

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Hexavalent Chromium	0.588	m mg/L	1.11	0.556	< 0.00659	105	80.1 - 118

⁴SPECIAL: Sodium Sulfate may have caused low LCS, LCSD, and MS recovery. Results may be biased low. •

⁵SPECIAL: Soduium Sulfate may have caused low LCS, LCSD, and MS recovery. Samples may be biased low. •

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	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Hexavalent Chromium	0.586	mg/L	1.11	0.556	< 0.00659	104	80.1 - 118	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63025 Date Analyzed: 2009-08-28 Analyzed By: ER
Prep Batch: 53796 QC Preparation: 2009-08-28 Prepared By: ER

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit GRO 0.920 < 0.07021.00 92 59.4 - 128 mg/L

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
GRO	0.932	mg/L	1	1.00	< 0.0702	93	59.4 - 128	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MS	MSD			Spike	MS	MSD	Rec .
Surrogate	Result	Result	Units	Dil.	${ m Amount}$	Rec .	Rec.	Limit
Trifluorotoluene (TFT)	0.0931	0.0870	mg/L	1	0.1	93	87	62.8 - 128
4-Bromofluorobenzene (4-BFB)	0.0836	0.0797	$_{ m mg/L}$	1	0.1	84	80	68.3 - 121

Matrix Spike (MS-1) Spiked Sample: 208427

QC Batch: 63042 Date Analyzed: 2009-08-28 Analyzed By: Prep Batch: 53812 QC Preparation: 2009-08-28 Prepared By:

	MS			${ m Spike}$	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec.}$	Limit
DRO	26.2	$\mathrm{mg/L}$	1	25.0	< 0.876	105	29.8 - 181

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	A mount	Result	Rec.	Limit	RPD	Limit
DRO	27.3	mg/L	1	25.0	< 0.876	109	29.8 - 181	4	20

	MS	MSD			Spike	MS	MSD	Rec .
Surrogate	Result	Result	Units	Dil.	${f Amount}$	$\mathrm{Rec.}$	$\mathrm{Rec.}$	Limit
n-Triacontane	8.56	8.80	m mg/L	1	10	86	88	34.4 - 185

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Analyzed By:

Prepared By:

TP

TP

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63072 Date Analyzed: Prep Batch: 53821

2009-08-31 QC Preparation: 2009-08-31

	MS			$_{ m Spike}$	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Total Mercury	0.000930	$_{ m mg/L}$	1	0.00100	< 0.0000329	93	80 - 116

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	Amount	Result	Rec .	Limit	RPD	Limit
Total Mercury	0.000930	mg/L	1	0.00100	< 0.0000329	93	80 - 116	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208427

2009-08-31 QC Batch: 63080 Date Analyzed: Analyzed By: AH Prep Batch: 53840QC Preparation: 2009-08-31 Prepared By: AH

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit
Total Cyanide	0.129	$_{ m mg/L}$	1	0.120	< 0.0110	108	66.3 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Total Cyanide	0.135	mg/L	1	0.120	< 0.0110	112	66.3 - 125	4	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208427

Analyzed By: AH QC Batch: 63111 Date Analyzed: 2009-08-31 Prep Batch: 53864QC Preparation: 2009-08-31 Prepared By: AH

	MS			Spike	Matrix		Rec .
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Ammonia-N	5.77	mg/L	1	5.00	0.392	108	57.2 - 133

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Ammonia-N	5.32	$\mathrm{mg/L}$	1	5.00	0.392	98	57.2 - 133	8	20

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Matrix Spike (MS-1) Spiked Sample: 208408

		MS			Spike	Matrix		${ m Rec.}$
Param		Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec}.$	Limit
HMX		1.60	$\mu { m g/L}$	1	2.50	< 0.123	64	10 - 164
RDX		1.86	$\mu { m g/L}$	1	2.50	< 0.298	74	10 - 147
1,3,5-Trinitrobenzene		1.64	$\mu { m g}/{ m L}$	1	2.50	< 0.339	66	10 - 187
1,3-Dinitrobenzene		1.82	$\mu { m g}/{ m L}$	1	2.50	< 0.389	73	10 - 155
Nitrobenzene		1.34	$\mu { m g}/{ m L}$	1	2.50	< 0.379	54	10 - 156
Tetryl		1.63	$\mu { m g}/{ m L}$	1	2.50	< 0.413	65	10 - 158
TNT		1.77	$\mu { m g}/{ m L}$	1	2.50	< 0.464	71	21 - 114
4-Amino-DNT	6	1.68	$\mu { m g/L}$	1	2.50	< 0.319	67	80 - 120
$2 ext{-Amino-DNT}$	7	1.77	$\mu { m g}/{ m L}$	1	2.50	< 0.391	71	80 - 120
2,6-DNT	8	1.60	$\mu { m g/L}$	1	2.50	< 0.323	64	80 - 120
2,4-DNT	9	1.80	$\mu { m g}/{ m L}$	1	2.50	< 0.366	72	80 - 120
2-NT	10	5.04	$\mu { m g}/{ m L}$	1	2.50	< 0.379	202	10 - 147
4-NT		1.40	$\mu { m g/L}$	1	2.50	< 0.398	56	10 - 161
3-NT		1.28	$\mu { m g}/{ m L}$	1	2.50	< 0.346	51	10 - 167

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

		MSD			Spike	Matrix		Rec .		RPD
Param		Result	Units	Dil.	A mount	Result	Rec .	Limit	RPD	Limit
HMX		1.66	$\mu \mathrm{g/L}$	1	2.50	< 0.123	66	10 - 164	4	20
RDX		1.76	$\mu { m g/L}$	1	2.50	< 0.298	70	10 - 147	6	20
1,3,5-Trinitrobenzene		1.92	$\mu { m g/L}$	1	2.50	< 0.339	77	10 - 187	16	20
1,3-Dinitrobenzene		1.94	$\mu { m g/L}$	1	2.50	< 0.389	78	10 - 155	6	20
${ m Nitrobenzene}$		1.39	$\mu { m g/L}$	1	2.50	< 0.379	56	10 - 156	4	20
Tetryl		1.56	$\mu { m g}/{ m L}$	1	2.50	< 0.413	62	10 - 158	4	20
TNT		1.83	$\mu { m g/L}$	1	2.50	< 0.464	73	21 - 114	3	20
4-Amino-DNT	11	1.71	$\mu { m g/L}$	1	2.50	< 0.319	68	80 - 120	2	20
2-Amino-DNT	12	1.98	$\mu { m g/L}$	1	2.50	< 0.391	79	80 - 120	11	20
2,6-DNT	13	1.70	$\mu \mathrm{g/L}$	1	2.50	< 0.323	68	80 - 120	6	20
2,4-DNT	14	1.88	$\mu { m g}/{ m L}$	1	2.50	< 0.366	75	80 - 120	4	20
2-NT	15	5.10	$\mu { m g}/{ m L}$	1	2.50	< 0.379	204	10 - 147	1	20
4-NT		1.36	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.398	54	10 - 161	3	20
3-NT		1.24	$\mu { m g}/{ m L}$	1	2.50	< 0.346	50	10 - 167	3	20

 $^{^6}$ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

⁷ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

⁸ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

⁹ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹⁰Matrix spike recovery out of control limits due to peak interference. Use LCS/LCSD to demonstrate analysis is under control.

¹¹ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹² Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.
¹³ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹⁴Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹⁵Matrix spike recovery out of control limits due to peak interference. Use LCS/LCSD to demonstrate analysis is under control.

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	MS	MSD			Spike	MS	MSD	Rec.
Surrogate	Result	Result	Units	Dil.	${f Amount}$	Rec.	Rec .	Limit
1 2-Dinitrobenzene	1 73	1.76	ug/L	1	2.5	69	70	10 - 222

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Total Silver 0.1370.125< 0.00111110 75 - 125 mg/L

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Silver	0.135	mg/L	1	0.125	< 0.00111	108	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Result Limit Param Units Dil. Amount Result Rec. Total Aluminum 1.26 1.00 0.065 120 75 - 125 mg/L

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Aluminum	1.24	mg/L	1	1.00	0.065	118	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit < 0.00448Total Arsenic 0.506mg/L0.500 101 75 - 125

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RR

KV

	MSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Arsenic	0.496	mg/L	1	0.500	< 0.00448	99	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By:

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Total Barium 1.00 1.00 0.008 99 75 - 125 mg/L1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD Spike Matrix Rec. RPD Param Result Units Dil. Amount Result Rec. Limit RPD Limit Total Barium 0.98820 mg/L1.00 0.008 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit 75 - 125 Total Beryllium 0.0254mg/L0.0250< 0.000450102

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSDRPD Spike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit RPD Limit 0.02470.0250< 0.00045020 Total Beryllium mg/L75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSMatrix Spike Rec. Param Result Units Dil. Amount Result Rec. Limit 0.227 0.250< 0.000303 Total Cadmium mg/L91 75 - 125

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KV

	MSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Total Cadmium	0.222	mg/L	1	0.250	< 0.000303	89	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

2009-09-03 QC Batch: 63221 Date Analyzed: Analyzed By: RR Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By:

MSMatrix Spike Rec. Param Result Units Dil. Amount Result Rec. Limit Total Cobalt 0.230 0.250 < 0.00082275 - 125 mg/L92 1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Cobalt	0.226	mg/L	1	0.250	< 0.000822	90	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: Date Analyzed: 2009-09-03 Analyzed By: RR 63221 Prep Batch: Prepared By: 53951 QC Preparation: 2009-09-03 KV

MSSpike MatrixRec. Param Units Result Dil. ${
m Amount}$ Result Rec. Limit Total Chromium 0.111 mg/L0.100 0.01596 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Chromium	0.108	$_{ m mg/L}$	1	0.100	0.015	93	75 - 125	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: QC Preparation: 2009-09-03 Prepared By: KV53951

MSMatrix Rec. Spike Param Result Units Dil. Amount Result Rec. Limit 0.144 0.125 < 0.000843 115 Total Copper mg/L75 - 125

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	MSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Total Copper	0.140	mg/L	1	0.125	< 0.000843	112	75 - 125	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSMatrix Rec. Spike Param Result Units Dil. Amount Result Rec. Limit Total Iron 0.4880.500 < 0.00087275 - 125 mg/L98 1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Iron	0.462	mg/L	1	0.500	< 0.000872	92	75 - 125	6	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Total Manganese 0.235mg/L0.250< 0.000305 94 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Manganese	0.231	mg/L	1	0.250	< 0.000305	92	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Rec. Matrix Param Result Units Dil. Amount Result Rec. Limit 0.5000.211Total Molybdenum 112 0.771mg/L75 - 125

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	MSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	$\overline{\mathrm{Amount}}$	Result	Rec .	Limit	RPD	Limit
Total Molybdenum	0.758	mg/L	1	0.500	0.211	109	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSMatrix Spike Rec. Param Result Units Dil. Amount Result Rec. Limit Total Nickel 0.230 0.250< 0.0012192 75 - 125 mg/L1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD Spike Matrix RPD Rec. Param Result Units Dil. Amount Result Rec. Limit RPD Limit Total Nickel 0.22520 mg/L0.250< 0.0012175 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike MatrixRec. Param Units Result Dil. ${
m Amount}$ Result Rec. Limit Total Phosphorous 0.493mg/L0.5000.00475 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD RPD Spike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit RPD Limit 0.50020 Total Phosphorous 0.485mg/L0.00475 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit 0.412 Total Lead mg/L0.500< 0.0032682 75 - 125

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	MSD			$_{ m Spike}$	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Lead	0.403	mg/L	1	0.500	< 0.00326	81	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSMatrix Spike Rec. Param Result Units Dil. Amount Result Rec. Limit Total Antimony 0.246 0.250< 0.0044075 - 125 mg/L98 1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD Spike Matrix RPD Rec. Param Result Units Dil. Amount Result Rec. Limit RPD Limit Total Antimony 0.24220 mg/L0.250< 0.0044075 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Dil. Result Units A mount Result Rec. Limit 75 - 125 Total Selenium 0.502mg/L0.500< 0.00508100

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD RPD Spike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit RPD Limit 0.500 20 Total Selenium 0.487mg/L< 0.00508 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Total Thallium 0.439mg/L0.500< 0.0048888 75 - 125

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	MSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	$\mathbf{A}\mathbf{mount}$	Result	Rec.	Limit	RPD	Limit
Total Thallium	0.437	mg/L	1	0.500	< 0.00488	87	75 - 125	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Total Vanadium 0.2650.250 0.017 99 75 - 125 mg/L1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD Spike Matrix Rec. RPD Param Result Units Dil. Amount Result Rec. Limit RPD Limit Total Vanadium 0.26020 mg/L0.2500.01797 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit 75 - 125 Total Zinc 0.258mg/L0.250< 0.000465103

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSDRPD Spike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit RPD Limit 0.253 0.25020 Total Zinc mg/L< 0.000465101 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208666

QC Batch: 63249 Date Analyzed: 2009-08-31 Analyzed By: SS Prep Batch: 53984 QC Preparation: 2009-08-31 Prepared By: SS

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit 276 250 110 90 - 110 Bromide mg/L50 < 9.60

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SS

	MSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Bromide	275	mg/L	50	250	< 9.60	110	90 - 110	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208666

QC Batch: 63249 Date Analyzed: 2009-08-31 Analyzed By: Prep Batch: 53984 QC Preparation: 2009-08-31 Prepared By:

MS Spike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit 2130 Chloride 50 1250 < 7.85170 90 - 110 mg/L

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD Spike Matrix Rec. RPD Param Result Units Dil. Amount Result Rec. Limit RPD Limit Chloride 2150 1250 90 - 110 20 mg/L50 < 7.85172

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208666

QC Batch: 63249 Date Analyzed: 2009-08-31 Analyzed By: SS Prep Batch: 53984 QC Preparation: 2009-08-31 Prepared By: SS

MSSpike Matrix Rec. Param Result Units Dil. ${
m Amount}$ Result Rec. Limit Fluoride 321 mg/L50 250 < 10.2128 90 - 110

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD RPD Spike Matrix Rec. Limit Param Result Units Dil. Amount Result Rec. RPD Limit 334 250 <10.2 134 20 Fluoride mg/L50 90 - 110

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208666

QC Batch: 63249 Date Analyzed: 2009-08-31 Analyzed By: SS Prep Batch: 53984 QC Preparation: 2009-08-31 Prepared By: SS

¹⁶Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹⁷ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹⁸Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹⁹ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

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		MS			Spike	Matrix		$\mathrm{Rec.}$
Param		Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Sulfate	20	5060	mg/L	50	1250	3460	128	90 - 110

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

		MSD			$_{ m Spike}$	Matrix		Rec .		RPD
Param		Result	Units	Dil.	A mount	Result	Rec.	Limit	RPD	Limit
Sulfate	21	5100	$\mathrm{mg/L}$	50	1250	3460	131	90 - 110	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63261 Date Analyzed: 2009-09-04 Analyzed By: MN Prep Batch: 53996 QC Preparation: 2009-09-02 Prepared By: MN

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Phenol	0.0194	$\mathrm{mg/L}$	1.02	0.0800	< 0.000519	24	10 - 66.5
2-Chlorophenol	0.0437	${ m mg/L}$	1.02	0.0800	< 0.000548	55	11.2 - 108
1,4-Dichlorobenzene (para)	0.0354	${ m mg/L}$	1.02	0.0800	< 0.000449	44	16 - 101
N-Nitrosodi-n-propylamine	0.0490	${ m mg/L}$	1.02	0.0800	< 0.000747	61	10 - 142
1,2,4-Trichlorobenzene	0.0389	${ m mg/L}$	1.02	0.0800	< 0.000412	49	18 - 108
Naphthalene	0.0420	${ m mg/L}$	1.02	0.0800	< 0.000499	52	20.2 - 114
4-Chloro-3-methylphenol	0.0558	${ m mg/L}$	1.02	0.0800	< 0.000532	70	21.5 - 125
${ m Acenapht hylene}$	0.0543	$\mathrm{mg/L}$	1.02	0.0800	< 0.000598	68	25.8 - 121
${ m Acenaphthene}$	0.0538	${ m mg/L}$	1.02	0.0800	< 0.000431	67	33.5 - 122
4-Nitrophenol	0.0218	${ m mg/L}$	1.02	0.0800	< 0.00189	27	10 - 125
2,4-Dinitrotoluene	0.0604	${ m mg/L}$	1.02	0.0800	< 0.000929	76	53 - 130
Fluorene	0.0572	${ m mg/L}$	1.02	0.0800	< 0.000661	72	44.6 - 117
Pentachlorophenol	0.0394	${ m mg/L}$	1.02	0.0800	< 0.000444	49	10 - 139
${ m Anthracene}$	0.0557	$\mathrm{mg/L}$	1.02	0.0800	< 0.000436	70	57.5 - 115
Phenanthrene	0.0558	$\mathrm{mg/L}$	1.02	0.0800	< 0.000559	70	55.5 - 118
Fluoranthene	0.0588	$\mathrm{mg/L}$	1.02	0.0800	< 0.000645	74	57 - 122
Pyrene	0.0524	${ m mg/L}$	1.02	0.0800	< 0.000737	66	58.5 - 130
$\operatorname{Benzo}(\operatorname{a})\operatorname{anthracene}$	0.0508	${ m mg/L}$	1.02	0.0800	< 0.000538	64	63.4 - 109
Chrysene	0.0568	${ m mg/L}$	1.02	0.0800	< 0.000651	71	54.7 - 114
Benzo(b)fluoranthene	0.0546	${ m mg/L}$	1.02	0.0800	< 0.000896	68	64.8 - 120
Benzo(k)fluoranthene	0.0721	${ m mg/L}$	1.02	0.0800	< 0.000862	90	70.3 - 114
Benzo(a)pyrene	0.0675	${ m mg/L}$	1.02	0.0800	< 0.00170	84	63.7 - 120
Indeno(1,2,3-cd)pyrene	0.0643	${ m mg/L}$	1.02	0.0800	< 0.000879	80	65.4 - 119
${ m Dibenzo(a,h)}$ anthracene	0.0634	$\mathrm{mg/L}$	1.02	0.0800	< 0.000825	79	68.7 - 117
$\operatorname{Benzo}(g,h,i)$ perylene	0.0664	$\mathrm{mg/L}$	1.02	0.0800	< 0.000968	83	57.2 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result. $continued \dots$

²⁰Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

²¹ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

$matrix\ spikes\ continued\ \dots$										
		MSD			Spike	Matrix		${ m Rec.}$		RPD
Param		Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit	RPD	Limit
		MSD			Spike	Matrix		Rec.		RPD
Param		Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Phenol		0.0182	mg/L	1.02	0.0800	< 0.000519	23	10 - 66.5	6	20
2-Chlorophenol		0.0132 0.0398	$\frac{\mathrm{mg}}{\mathrm{L}}$	1.02	0.0800	< 0.000513	$\frac{23}{50}$	11.2 - 108	9	$\frac{20}{20}$
1,4-Dichlorobenzene (para)		0.0398 0.0317	$\frac{\mathrm{mg}}{\mathrm{L}}$	1.02 1.02	0.0800	< 0.000348	40	16 - 101	9 11	$\frac{20}{20}$
N-Nitrosodi-n-propylamine		0.0317 0.0465	$\frac{\mathrm{mg}}{\mathrm{L}}$	1.02 1.02	0.0800	< 0.000449	58	10 - 101	5	$\frac{20}{20}$
1,2,4-Trichlorobenzene		0.0403 0.0344	$\frac{\mathrm{mg}}{\mathrm{L}}$	1.02 1.02	0.0800	< 0.000412	43	18 - 108	12	$\frac{20}{20}$
Naphthalene		0.0344 0.0363	$\frac{\mathrm{mg}}{\mathrm{L}}$	1.02 1.02	0.0800	< 0.000412	$\frac{45}{45}$	20.2 - 114	15	$\frac{20}{20}$
4-Chloro-3-methylphenol		0.0303 0.0496	$\frac{\mathrm{mg}}{\mathrm{L}}$	1.02 1.02	0.0800	< 0.000499	62	21.5 - 125	$\frac{13}{12}$	$\frac{20}{20}$
Acenaphthylene		0.0430 0.0479	$\frac{\mathrm{mg}}{\mathrm{L}}$	1.02	0.0800	< 0.000598	60	25.8 - 121	$\frac{12}{12}$	$\frac{20}{20}$
Acenaphthene		0.0479	$\frac{\mathrm{mg}}{\mathrm{L}}$	1.02	0.0800	< 0.000333	58	33.5 - 122	$\frac{12}{14}$	$\frac{20}{20}$
4-Nitrophenol		0.0408 0.0200		1.02	0.0800	< 0.000431	$\frac{36}{25}$	10 - 125	9	$\frac{20}{20}$
2,4-Dinitrotoluene		0.0200 0.0526	${ m mg/L} \ { m mg/L}$	1.02 1.02	0.0800	< 0.00139	66	53 - 130	9 14	$\frac{20}{20}$
Fluorene		0.0320 0.0497		1.02 1.02	0.0800	< 0.000929	62	44.6 - 117	14	$\frac{20}{20}$
Pentachlorophenol		0.0497 0.0358	mg/L	1.02 1.02	0.0800	< 0.000444	$\frac{62}{45}$	10 - 139	10	$\frac{20}{20}$
Anthracene		0.0508	$\frac{\text{mg/L}}{\text{mg/L}}$	1.02 1.02	0.0800	< 0.000444	63	57.5 - 115	10	$\frac{20}{20}$
Phenanthrene		0.0303 0.0492	$\frac{\text{mg/L}}{\text{mg/L}}$	1.02 1.02	0.0800	< 0.000450 < 0.000559	62	55.5 - 118	$\frac{10}{13}$	$\frac{20}{20}$
Fluoranthene		0.0492 0.0537	$\frac{\text{mg/L}}{\text{mg/L}}$	1.02 1.02	0.0800	< 0.000645	67	55.5 - 116 57 - 122	13 9	$\frac{20}{20}$
	22		$\frac{\text{mg/L}}{\text{mg/L}}$	1.02 1.02	0.0800		57	58.5 - 130	$\frac{9}{14}$	$\frac{20}{20}$
Pyrene Panya (a) anthrasana	23	0.0457 0.0446	$\frac{\text{mg/L}}{\text{mg/L}}$	1.02 1.02	0.0800	< 0.000737 < 0.000538	56	63.4 - 109	$\frac{14}{13}$	$\frac{20}{20}$
Benzo(a)anthracene			mg/L				50 61			$\frac{20}{20}$
Chrysene		0.0485	mg/L	1.02	0.0800	< 0.000651		54.7 - 114	16	
Benzo(b)fluoranthene	24	0.0528	mg/L	1.02	0.0800	< 0.000896	66	64.8 - 120	3	20
Benzo(k)fluoranthene		0.0573	mg/L	1.02	0.0800	< 0.000862	72	70.3 - 114	23	20
Benzo(a) pyrene		0.0589	mg/L	1.02	0.0800	< 0.00170	74	63.7 - 120	14	20
Indeno(1,2,3-cd)pyrene	25	0.0540	mg/L	1.02	0.0800	< 0.000879	68	65.4 - 119	17	20
Dibenzo(a,h)anthracene	20	0.0536	mg/L	1.02	0.0800	< 0.000825	67	68.7 - 117	17	20
Benzo(g,h,i)perylene		0.0557	${ m mg/L}$	1.02	0.0800	< 0.000968	70	57.2 - 125	18	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MS	MSD			Spike	MS	MSD	$\mathrm{Rec}.$
$\operatorname{Surrogate}$	Result	Result	Units	Dil.	${ m Amount}$	Rec.	Rec .	Limit
2-Fluorophenol	0.0262	0.0254	m mg/L	1.02	0.08	33	32	10 - 53.1
${ m Phenol-d5}$	0.0194	0.0183	${ m mg/L}$	1.02	0.08	24	23	10 - 36.9
${ m Nitrobenzene-d5}$	0.0450	0.0403	${ m mg/L}$	1.02	0.08	56	50	23.8 - 108
2-Fluorobiphenyl	0.0445	0.0387	${ m mg/L}$	1.02	0.08	56	48	15.9 - 127
2,4,6-Tribromophenol	0.0653	0.0583	${ m mg/L}$	1.02	0.08	82	73	10 - 123
Terphenyl-d14	0.0522	0.0450	${ m mg/L}$	1.02	0.08	65	56	17.2 - 160

Matrix Spike (MS-1) Spiked Sample: 208665

QC Batch: 63291 Date Analyzed: 2009-09-02 Analyzed By: AH Prep Batch: 54023 QC Preparation: 2009-09-02Prepared By: AH

²²Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

²³ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

 $^{^{24}\,\}mathrm{MS/MSD}$ RPD out of RPD Limits. Use LCS/LCSD to demonstrate analysis is under control.

²⁵Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

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		MS			Spike	Matrix		Rec.
Param		Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Kjeldahl Nitrogen - N	26	33.9	mg/L	1	50.0	4.48	59	61.2 - 118

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Total Kjeldahl Nitrogen - N	38.1	$\mathrm{mg/L}$	1	50.0	4.48	67	61.2 - 118	12	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63340 Date Analyzed: 2009-09-08 Analyzed By: KV Prep Batch: 54065 QC Preparation: 2009-09-08 Prepared By: KV

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Organic Carbon	51.2	mg/L	1	50.0	1.66	99	66.9 - 121

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Organic Carbon	48.4	mg/L	1	50.0	1.66	93	66.9 - 121	6	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Total Calcium	574	mg/L	1	50.0	531	86	75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Calcium	576	mg/L	1	50.0	531	90	75 - 125	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

²⁶ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

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	MS			Spike	Matrix		Rec.
Param	Result	Units	Dil.	$\overline{ ext{Amount}}$	Result	Rec .	Limit
Total Potassium	150	mg/L	1	50.0	99.9	100	75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			$_{ m Spike}$	Matrix		Rec .		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Potassium	148	mg/L	1	50.0	99.9	96	75 - 125	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

	MS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${ m Amount}$	Result	$\mathrm{Rec}.$	Limit
Total Magnesium	471	$\mathrm{mg/L}$	1	50.0	425	92	75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Magnesium	472	mg/L	1	50.0	425	94	75 - 125	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

	MS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Total Sodium	768	mg/L	1	50.0	717	102	75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec.}$		RPD
Param	Result	Units	Dil .	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Sodium	776	mg/L	1	50.0	717	118	75 - 125	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63461 Date Analyzed: 2009-09-11 Analyzed By: RR
Prep Batch: 54153 QC Preparation: 2009-09-11 Prepared By: KV

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	MS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Dissolved Chromium	0.110	mg/L	1	0.100	0.014	96	75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit	RPD	Limit
Dissolved Chromium	0.112	${ m mg/L}$	1	0.100	0.014	98	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208427

QC Batch: 63485 Date Analyzed: 2009-09-11 Analyzed By: KV Prep Batch: 54188 QC Preparation: 2009-09-11 Prepared By: KV

		MS			Spike	Matrix		${ m Rec.}$
Param		Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Nitrate and Nitrite as N	27	19.2	$_{ m mg/L}$	200	0.200	14.5	2350	80 - 120

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

		MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param		Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Nitrate and Nitrite as N	28	18.7	mg/L	200	0.200	14.5	9350	80 - 120	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Standard (CCV-1)

QC Batch:	63015	Date Analyzed:	2009-08-27	Analyzed By: MD
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			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Hexavalent Chromium		m mg/L	0.500	0.501	100	90 - 110	2009-08-27

Standard (CCV-2)

QC Batch: 63015 Date Analyzed: 2009-08-27 Analyzed By: MD

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	${ m Units}$	Conc .	Conc .	Recovery	Limits	${\bf Analyzed}$
Hexavalent Chromium		${ m mg/L}$	0.500	0.497	99	90 - 110	2009-08-27

²⁷ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

²⁸Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

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Standard	(CCV-1)
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QC Batch: 63025 Date Analyzed: 2009-08-28 Analyzed By: ER

CCVsCCVs CCVs Percent TrueFound Percent Recovery Date Param Flag Units Conc. Conc. Recovery Limits Analyzed $\overline{\text{GRO}}$ 80 - 120 2009-08-28 mg/L1.00 0.95796

Standard (CCV-2)

QC Batch: 63025 Date Analyzed: 2009-08-28 Analyzed By: ER

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
GRO		$_{ m mg/L}$	1.00	0.928	93	80 - 120	2009-08-28

Standard (CCV-1)

QC Batch: 63042 Date Analyzed: 2009-08-28 Analyzed By:

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
DRO		$\mathrm{mg/L}$	250	300	120	80 - 120	2009-08-28

Standard (CCV-2)

QC Batch: 63042 Date Analyzed: 2009-08-28 Analyzed By:

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
DRO		$\mathrm{mg/L}$	250	290	116	80 - 120	2009-08-28

Standard (ICV-1)

Analyzed By: TP QC Batch: 63072 Date Analyzed: 2009-08-31

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Mercury		$\mathrm{mg/L}$	0.00100	0.00102	102	90 - 110	2009-08-31

Standard (CCV-1)

QC Batch: 63072Date Analyzed: 2009-08-31 Analyzed By: TP

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HELSTF Chromate Spill Groundwater

			CCVs	CCVs	CCVs	Percent	D-4-
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	${ m Units}$	Conc .	$\operatorname{Conc.}$	Recovery	Limits	\mathbf{A} nalyzed
Total Mercury		m mg/L	0.00100	0.000950	95	90 - 110	2009-08-31

Standard (ICV-1)

QC Batch: 63080

Date Analyzed: 2009-08-31

Analyzed By: AH

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			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Cyanide		$\mathrm{mg/L}$	0.120	0.117	98	85 - 115	2009-08-31

Standard (CCV-1)

QC Batch: 63080

Date Analyzed: 2009-08-31

Analyzed By: AH

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Cyanide		m mg/L	0.120	0.120	100	85 - 115	2009-08-31

Standard (ICV-1)

QC Batch: 63082

Date Analyzed: 2009-08-27

Analyzed By: JG

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
$\overline{\mathrm{pH}}$		s.u.	7.00	7.00	100	98 - 102	2009-08-27

Standard (CCV-1)

QC Batch: 63082

Date Analyzed: 2009-08-27

Analyzed By: JG

Analyzed By: AH

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
pН		s.u.	7.00	7.00	100	98 - 102	2009-08-27

Standard (ICV-1)

QC Batch: 63111

Work Order: 9082815

HELSTF Chromate Spill Groundwater

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	Units	Conc.	Conc .	Recovery	Limits	${ m Analyzed}$
Ammonia-N		$_{ m mg/L}$	5.00	4.82	96	85 - 115	2009-08-31

Standard (CCV-1)

QC Batch: 63111

Date Analyzed: 2009-08-31

Analyzed By: AH

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			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Ammonia-N		$\mathrm{mg/L}$	5.00	4.79	96	85 - 115	2009-08-31

Standard (ICV-1)

QC Batch: 63142

Date Analyzed: 2009-09-01

Analyzed By: DS

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	$_{ m Units}$	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
HMX		$\mu { m g/L}$	500	533	107	85 - 115	2009-09-01
RDX		$\mu { m g}/{ m L}$	500	487	97	85 - 115	2009-09-01
1,3,5-Trinitrobenzene		$\mu { m g}/{ m L}$	500	503	101	85 - 115	2009-09-01
1,3-Dinitrobenzene		$\mu { m g}/{ m L}$	500	491	98	85 - 115	2009-09-01
${ m Nitrobenzene}$		$\mu { m g}/{ m L}$	500	515	103	85 - 115	2009-09-01
Tetryl		$\mu { m g}/{ m L}$	500	501	100	85 - 115	2009-09-01
TNT		$\mu \mathrm{g}/\mathrm{L}$	500	498	100	85 - 115	2009-09-01
4-Amino-DNT		$\mu { m g}/{ m L}$	500	497	99	85 - 115	2009-09-01
2-Amino-DNT		$\mu { m g}/{ m L}$	500	529	106	85 - 115	2009-09-01
2,6-DNT		$\mu \mathrm{g}/\mathrm{L}$	500	440	88	85 - 115	2009-09-01
2,4-DNT		$\mu \mathrm{g}/\mathrm{L}$	500	525	105	85 - 115	2009-09-01
2-NT		$\mu { m g}/{ m L}$	500	480	96	85 - 115	2009-09-01
4-NT		$\mu \mathrm{g}/\mathrm{L}$	500	493	99	85 - 115	2009-09-01
3-NT		$\mu { m g}/{ m L}$	500	506	101	85 - 115	2009-09-01

					Spike	$\operatorname{Percent}$	Recovery
$\mathbf{Surrogate}$	Flag	Result	Units	Dilution	${ m Amount}$	Recovery	Limit
1,2-Dinitrobenzene		494	$\mu \mathrm{g}/\mathrm{L}$	1	500	99	85 - 115

Standard (CCV-1)

QC Batch: 63142

Date Analyzed: 2009-09-01

Analyzed By: DS

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
HMX		$\mu { m g/L}$	500	502	100	85 - 115	2009-09-01
RDX		$\mu { m g}/{ m L}$	500	499	100	85 - 115	2009-09-01

 $continued \dots$

at and and	ann timu ad	

source constituee			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	$_{ m Units}$	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
1,3,5-Trinitrobenzene		$\mu \mathrm{g/L}$	500	500	100	85 - 115	2009-09-01
1,3-Dinitrobenzene		$\mu { m g}/{ m L}$	500	507	101	85 - 115	2009-09-01
Nitrobenzene		$\mu { m g}/{ m L}$	500	542	108	85 - 115	2009-09-01
Tetryl		$\mu { m g}/{ m L}$	500	495	99	85 - 115	2009-09-01
TNT		$\mu { m g}/{ m L}$	500	534	107	85 - 115	2009-09-01
4-Amino-DNT		$\mu { m g}/{ m L}$	500	553	111	85 - 115	2009-09-01
2-Amino-DNT		$\mu { m g}/{ m L}$	500	538	108	85 - 115	2009-09-01
2,6-DNT		$\mu { m g}/{ m L}$	500	461	92	85 - 115	2009-09-01
2,4-DNT		$\mu { m g}/{ m L}$	500	557	111	85 - 115	2009-09-01
2-NT		$\mu { m g}/{ m L}$	500	516	103	85 - 115	2009-09-01
4-NT		$\mu { m g}/{ m L}$	500	515	103	85 - 115	2009-09-01
3-NT		$\mu { m g}/{ m L}$	500	550	110	85 - 115	2009-09-01

					Spike	$\operatorname{Percent}$	$\operatorname{Recovery}$
$\mathbf{Surrogate}$	Flag	Result	Units	Dilution	${f Amount}$	Recovery	Limit
1,2-Dinitrobenzene		515	$\mu \mathrm{g/L}$	1	500	103	85 - 115

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Page Number: 65 of 79

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Total Silver		$_{ m mg/L}$	0.250	0.253	101	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Aluminum		$_{ m mg/L}$	1.00	1.01	101	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
Total Arsenic		$_{ m mg/L}$	2.00	2.02	101	90 - 110	2009-09-03

Work Order: 9082815 HELSTF Chromate Spill Groundwater

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Standard (1	ICV-1)
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QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Barium		$_{ m mg/L}$	1.00	1.02	102	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Beryllium		mg/L	1.00	1.02	102	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Total Cadmium		mg/L	1.00	1.04	104	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
Total Cobalt		m mg/L	1.00	0.992	99	90 - 110	2009-09-03

Standard (ICV-1)

Analyzed By: RR QC Batch: 63221 Date Analyzed: 2009-09-03

			$\mathrm{CC}\mathrm{Vs}$	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Total Chromium		mg/L	1.00	1.04	104	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221Date Analyzed: 2009-09-03 Analyzed By: RR

Work Order: 9082815

HELSTF Chromate Spill Groundwater

			CCVs True	CCVs Found	CCVs Percent	Percent Recovery	Date
Param	Flag	Units	Conc.	Conc.	Recovery	Limits	$\begin{array}{c} \text{Analyzed} \end{array}$
Total Copper		$\mathrm{mg/L}$	1.00	1.05	105	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Page Number: 67 of 79

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	$\operatorname{Recovery}$	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Iron		$_{ m mg/L}$	1.00	1.04	104	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Manganese		$_{ m mg/L}$	1.00	1.01	101	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Molybdenum		$_{ m mg/L}$	1.00	0.992	99	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Nickel		mg/L	1.00	0.999	100	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Work Order: 9082815 HELSTF Chromate Spill Groundwater

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${\bf Analyzed}$
Total Phosphorous		$\mathrm{mg/L}$	5.00	4.93	99	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Page Number: 68 of 79

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Lead		$\mathrm{mg/L}$	2.00	2.06	103	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Antimony		$\mathrm{mg/L}$	2.00	2.04	102	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Total Selenium		$_{ m mg/L}$	1.00	1.02	102	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
Total Thallium		$\mathrm{mg/L}$	5.00	5.10	102	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Work Order: 9082815 HELSTF Chromate Spill Groundwater

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Total Vanadium		mg/L	1.00	1.04	104	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

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			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	$\operatorname{Recovery}$	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Total Zinc		$_{ m mg/L}$	1.00	1.08	108	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
Total Silver		$_{ m mg/L}$	0.125	0.129	103	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Aluminum		$_{ m mg/L}$	1.00	0.993	99	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Arsenic		$_{ m mg/L}$	1.00	0.981	98	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

HELSTF Chromate Spill Groundwater

Work Order: 9082815	Page Number: 70 of 79
TOTE Charmete Caill Casuadaneter	

			CCVs True	CCVs Found	${ m CCVs} \ { m Percent}$	Percent Recovery	Date
Param	Flag	Units	Conc.	Conc.	Recovery	$rac{ ext{Limits}}{ ext{Limits}}$	$\begin{array}{c} \text{Date} \\ \text{Analyzed} \end{array}$
Total Barium		mg/L	1.00	1.04	104	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Beryllium		mg/L	1.00	1.00	100	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221 Analyzed By: RR Date Analyzed: 2009-09-03

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Cadmium		mg/L	1.00	1.01	101	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Cobalt		$\mathrm{mg/L}$	1.00	0.989	99	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Chromium		mg/L	1.00	1.02	102	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

mg/L

Work Order: 9082815 HELSTF Chromate Spill Groundwater

1.04

		CCVs	CCVs	CCVs	Percent	
		True	Found	$\operatorname{Percent}$	Recovery	Date
Flag	Units	Conc	Conc	Recovery	Limits	Analyzed

104

Standard (CCV-1)

QC Batch: 63221

Param

Total Copper

Date Analyzed: 2009-09-03

1.00

Analyzed By: RR

2009-09-03

90 - 110

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			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Total Iron		$\mathrm{mg/L}$	1.00	1.02	102	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Manganese		$\mathrm{mg/L}$	1.00	0.996	100	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Molybdenum		$_{ m mg/L}$	1.00	0.985	98	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Nickel		$_{ m mg/L}$	1.00	0.973	97	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Work Order: 9082815 HELSTF Chromate Spill Groundwater

Work Order: 9082815	Page Number: 72 of 79
ELSTE Chromate Spill Groundwater	

			$_{ m CCVs}$	CCVs	$_{ m CCVs}$	Percent	D .
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	${ m Units}$	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Phosphorous		${ m mg/L}$	5.00	4.84	97	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Lead		$\mathrm{mg/L}$	1.00	0.966	97	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Antimony		$\mathrm{mg/L}$	1.00	1.00	100	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Total Selenium		$_{ m mg/L}$	1.00	0.983	98	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
Total Thallium		$\mathrm{mg/L}$	1.00	1.00	100	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

				1			
			$rac{ ext{CCVs}}{ ext{True}}$	$\begin{array}{c} { m CCVs} \\ { m Found} \end{array}$	$rac{ ext{CCVs}}{ ext{Percent}}$	Percent Recovery	Date
Param	Flag	Units	Conc.	Conc.	Recovery	Limits	$rac{ m Date}{ m Analyzed}$
Total Vanadium		$\mathrm{mg/L}$	1.00	1.04	104	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Page Number: 73 of 79

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Total Zinc		mg/L	1.00	1.03	103	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63249

Date Analyzed: 2009-08-31

Analyzed By: SS

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Bromide		$\mathrm{mg/L}$	5.00	5.05	101	90 - 110	2009-08-31

Standard (CCV-1)

QC Batch: 63249

Date Analyzed: 2009-08-31

Analyzed By: SS

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Chloride		mg/L	25.0	23.7	95	90 - 110	2009-08-31

Standard (CCV-1)

QC Batch: 63249

Date Analyzed: 2009-08-31

Analyzed By: SS

Analyzed By: SS

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
Fluoride		$_{ m mg/L}$	5.00	5.33	107	90 - 110	2009-08-31

Standard (CCV-1)

QC Batch: 63249

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			пегот	r Cinomate 51	oni Groundwater		
Param	Floor	Units	$\begin{array}{c} { m CCVs} \\ { m True} \\ { m Conc.} \end{array}$	$\begin{array}{c} {\rm CCVs} \\ {\rm Found} \\ {\rm Conc.} \end{array}$	CCVs Percent	Percent Recovery Limits	Date
Sulfate	Flag	mg/L	25.0	24.7	Recovery 99	90 - 110	Analyzed 2009-08-31
<u>sunate</u>		mg/L	29.0	24.1	99	90 - 110	2009-08-31
Standard ((CCV-2)						
QC Batch:	63249		Date An	alyzed: 2009-	08-31	Ana	lyzed By: SS
			CCVs	CCVs	CCVs	Percent	
_			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc.	Conc.	Recovery	Limits	Analyzed
Bromide		m mg/L	5.00	5.04	101	90 - 110	2009-08-31
Standard ((CCV-2)						
QC Batch:	63249		Date An	alyzed: 2009-	08-31	Ana	lyzed By: SS
			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc.	Conc.	Recovery	Limits	Analyzed
Chloride		mg/L	25.0	23.9	96	90 - 110	2009-08-31
Standard ((CCV-2)						
QC Batch:	63249		Date An	alyzed: 2009-	08-31	Ana	lyzed By: SS
			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Fluoride		mg/L	5.00	5.38	108	90 - 110	2009-08-31
Standard ((CCV-2)						
QC Batch:	63249		Date An	alyzed: 2009-	08-31	Ana	lyzed By: SS
			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Sulfate	-	$\mathrm{mg/L}$	25.0	24.3	97	90 - 110	2009-08-31

Standard (CCV-1)

QC Batch: 63261 Date Analyzed: 2009-09-04 Analyzed By: MN

Diphenylamine

Fluoranthene

Benzo(a)pyrene

Pentachlorophenol

Di-n-octylphthalate

Work Order: 9082815 **HELSTF** Chromate Spill Groundwater Page Number: 75 of 79

2009 - 09 - 04

2009-09-04

2009-09-04

2009-09-04

2009-09-04

IIII om om oo ppin ground weer									
			CCVs	CCVs	CCVs	Percent			
			True	Found	Percent	Recovery	Date		
Param	Flag	Units	Conc.	$\operatorname{Conc.}$	Recovery	Limits	Analyzed		
Phenol		$_{ m mg/L}$	60.0	55.1	92	80 - 120	2009-09-04		
1,4-Dichlorobenzene (para)		${ m mg/L}$	60.0	59.4	99	80 - 120	2009-09-04		
2-Nitrophenol		m mg/L	60.0	65.0	108	80 - 120	2009-09-04		
2,4-Dichlorophenol		m mg/L	60.0	67.4	112	80 - 120	2009-09-04		
${ m Hexachlorobutadiene}$		m mg/L	60.0	61.7	103	80 - 120	2009-09-04		
4-Chloro-3-methylphenol	29	$_{ m mg/L}$	60.0	73.5	122	80 - 120	2009-09-04		
2,4,6-Trichlorophenol		m mg/L	60.0	63.5	106	80 - 120	2009-09-04		
${f Acenaphthene}$		mg/L	60.0	60.2	100	80 - 120	2009-09-04		

60.4

58.4

59.8

63.6

62.7

101

97

100

106

104

80 - 120

80 - 120

80 - 120

80 - 120

80 - 120

Surrogate	Flag	Result	Units	Dilution	$egin{aligned} ext{Spike} \ ext{Amount} \end{aligned}$	Percent Recovery	$egin{array}{c} ext{Recovery} \ ext{Limit} \end{array}$
2-Fluorophenol		61.7	$\mathrm{mg/L}$	1	60.0	103	80 - 120
${ m Phenol-d5}$		56.5	$\mathrm{mg/L}$	1	60.0	94	80 - 120
${ m Nitrobenzene-d5}$		58.2	$\mathrm{mg/L}$	1	60.0	97	80 - 120
2-Fluorobiphenyl		58.2	${ m mg/L}$	1	60.0	97	80 - 120
2,4,6-Tribromophenol		70.7	${ m mg/L}$	1	60.0	118	80 - 120
Terphenyl-d14		57.7	$\mathrm{mg/L}$	1	60.0	96	80 - 120

60.0

60.0

60.0

60.0

60.0

mg/L

mg/L

mg/L

mg/L

mg/L

Standard (ICV-1)

QC Batch: 63272 Date Analyzed: 2009-09-01 Analyzed By: MD

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc.	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Dissolved Solids		mg/L	1000	980	98	90 - 110	2009-09-01

Standard (CCV-1)

QC Batch: 63272 Date Analyzed: 2009-09-01 Analyzed By: MD

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc.	Conc .	Recovery	Limits	${\bf Analyzed}$
Total Dissolved Solids		mg/L	1000	1010	101	90 - 110	2009-09-01

Standard (ICV-1)

QC Batch: 63291 Date Analyzed: 2009-09-02 Analyzed By: AH

²⁹Control analyte out of CCV control limits. Results biased high. •

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
Total Kjeldahl Nitrogen - N		m mg/L	5.00	4.96	99	85 - 115	2009-09-02

Standard (CCV-1)

QC Batch: 63291

Date Analyzed: 2009-09-02

Analyzed By: AH

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			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc.	Conc .	Recovery	Limits	${ m Analyzed}$
Total Kjeldahl Nitrogen - N		m mg/L	5.00	4.87	97	85 - 115	2009-09-02

Standard (CCV-1)

QC Batch: 63340

Date Analyzed: 2009-09-08

Analyzed By: KV

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Organic Carbon		${ m mg/L}$	50.0	49.3	99	80 - 120	2009-09-08

Standard (CCV-2)

QC Batch: 63340

Date Analyzed: 2009-09-08

Analyzed By: KV

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Organic Carbon		${ m mg/L}$	50.0	49.8	100	80 - 120	2009-09-08

Standard (ICV-1)

QC Batch: 63355

Date Analyzed: 2009-09-09

Analyzed By: RR

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Calcium		$\mathrm{mg/L}$	50.0	51.5	103	90 - 110	2009-09-09

Standard (ICV-1)

QC Batch: 63355

Work Order: 9082815 HELSTF Chromate Spill Groundwater

-							
			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${\bf Analyzed}$
Total Potassium		$_{ m mg/L}$	50.0	50.8	102	90 - 110	2009-09-09

Standard (ICV-1)

QC Batch: 63355

Date Analyzed: 2009-09-09

Analyzed By: RR

Page Number: 77 of 79

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Magnesium		$\mathrm{mg/L}$	50.0	51.0	102	90 - 110	2009-09-09

Standard (ICV-1)

QC Batch: 63355

Date Analyzed: 2009-09-09

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Sodium		m mg/L	50.0	50.6	101	90 - 110	2009-09-09

Standard (CCV-1)

QC Batch: 63355

Date Analyzed: 2009-09-09

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${f Analyzed}$
Total Calcium		mg/L	50.0	51.1	102	90 - 110	2009-09-09

Standard (CCV-1)

QC Batch: 63355

Date Analyzed: 2009-09-09

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
Total Potassium		$\mathrm{mg/L}$	50.0	51.1	102	90 - 110	2009-09-09

Standard (CCV-1)

QC Batch: 63355

Date Analyzed: 2009-09-09

Analyzed By: RR

 ${\it Work~Order:~9082815} \\ {\it HELSTF~Chromate~Spill~Groundwater} \\$

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${\bf Analyzed}$
Total Magnesium		mg/L	50.0	50.4	101	90 - 110	2009-09-09

Standard (CCV-1)

QC Batch: 63355

Date Analyzed: 2009-09-09

Analyzed By: RR

Page Number: 78 of 79

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Total Sodium		$_{ m mg/L}$	50.0	50.1	100	90 - 110	2009-09-09

Standard (ICV-1)

QC Batch: 63379

Date Analyzed: 2009-09-08

Analyzed By: JG

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc.	Conc .	Recovery	Limits	${\bf Analyzed}$
Hydroxide Alkalinity		mg/L as CaCo3	0.00	<1.00		90 - 110	2009-09-08
Carbonate Alkalinity		mg/L as $CaCo3$	0.00	240		90 - 110	2009-09-08
Bicarbonate Alkalinity		mg/L as $CaCo3$	0.00	8.00		90 - 110	2009-09-08
Total Alkalinity		mg/L as $CaCo3$	250	248	99	90 - 110	2009-09-08

Standard (CCV-1)

QC Batch: 63379

Date Analyzed: 2009-09-08

Analyzed By: JG

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Hydroxide Alkalinity		mg/L as CaCo3	0.00	<1.00		90 - 110	2009-09-08
Carbonate Alkalinity		mg/L as $CaCo3$	0.00	240		90 - 110	2009-09-08
Bicarbonate Alkalinity		mg/L as $CaCo3$	0.00	10.0		90 - 110	2009-09-08
Total Alkalinity		mg/L as $CaCo3$	250	250	100	90 - 110	2009-09-08

Standard (ICV-1)

QC Batch: 63461

Date Analyzed: 2009-09-11

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	$_{ m Units}$	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Dissolved Chromium		mg/L	1.00	1.03	103	90 - 110	2009-09-11

Report Date: September 11, 2009 Work Order: 9082815 Page Number: 79 of 79

HELSTF Chromate Spill Groundwater

Standard (CCV-1)

QC Batch: 63461 Date Analyzed: 2009-09-11 Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Dissolved Chromium		mg/L	1.00	1.02	102	90 - 110	2009-09-11

Standard (ICV-1)

QC Batch: 63485 Date Analyzed: 2009-09-11 Analyzed By: KV

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Nitrate and Nitrite as N		mg/L	0.200	0.188	94	85 - 115	2009-09-11

Standard (CCV-1)

QC Batch: 63485 Date Analyzed: 2009-09-11 Analyzed By: KV

			$rac{ ext{CCVs}}{ ext{True}}$	$\begin{array}{c} { m CCVs} \\ { m Found} \end{array}$	$rac{ ext{CCVs}}{ ext{Percent}}$	Percent Recovery	Date
Param	Flag	Units	Conc.	$\operatorname{Conc.}$	Recovery	Limits	${\bf Analyzed}$
Nitrate and Nitrite as N		mg/L	0.200	0.201	100	85 - 115	2009-09-11

9082815

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